

V. Balasubramanian

Trying 3106016892...Open

Welcome to STN International! Enter x:x
LOGINID:sssptal611bxv
PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

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FILE 'HOME' ENTERED AT 15:32:00 ON 05 OCT 2000

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.15	0.15

FILE 'REGISTRY' ENTERED AT 15:32:06 ON 05 OCT 2000
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STRUCTURE FILE UPDATES: 4 OCT 2000 HIGHEST RN 292819-21-7
DICTIONARY FILE UPDATES: 4 OCT 2000 HIGHEST RN 292819-21-7

09/825,925

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TSCA INFORMATION NOW CURRENT THROUGH JANUARY 11, 2000

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT
for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>

Uploading c:\stnexp4\queries\09543703.str

L1 STRUCTURE UPLOADED

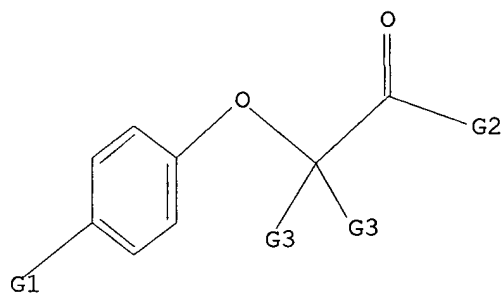
=> que L1

L2 QUE L1

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 C,H,O,N

G2 N,OH

G3 H,Me,Ak

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss sam

SAMPLE SEARCH INITIATED 15:32:42 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 11130 TO ITERATE

9.0% PROCESSED 1000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

50 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

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PROJECTED ITERATIONS: 216297 TO 228903
PROJECTED ANSWERS: 42770 TO 48496

L3 50 SEA SSS SAM L1

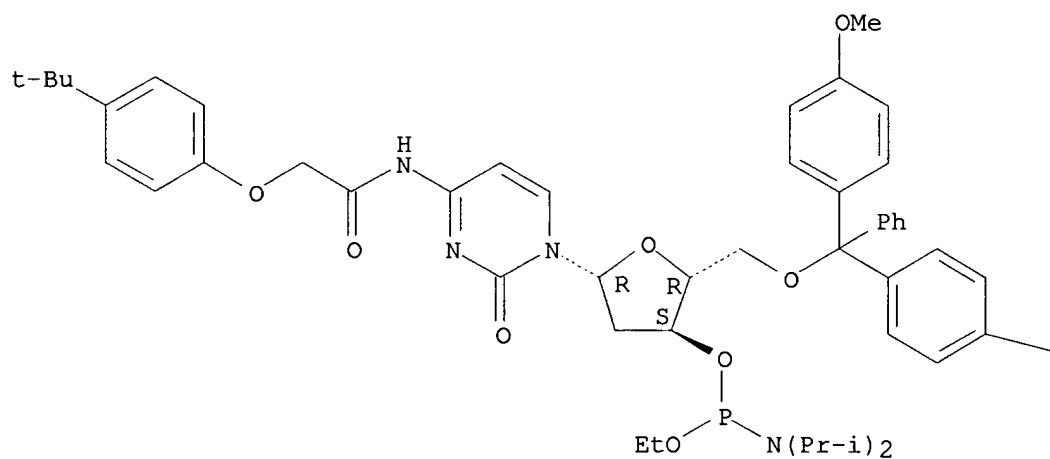
=> d scan 1-50

'1-50' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Cytidine, 5'-O-[bis(4-methoxyphenyl)phenylmethyl]-2'-deoxy-N-[[4-(1,1-dimethylethyl)phenoxy]acetyl]-, 3'-[ethyl bis(1-methylethyl)phosphoramidite] (9CI)
MF C50 H63 N4 O9 P

Absolute stereochemistry.

PAGE 1-A



PAGE 1-B

— OMe

The following are valid formats:

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Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN
SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

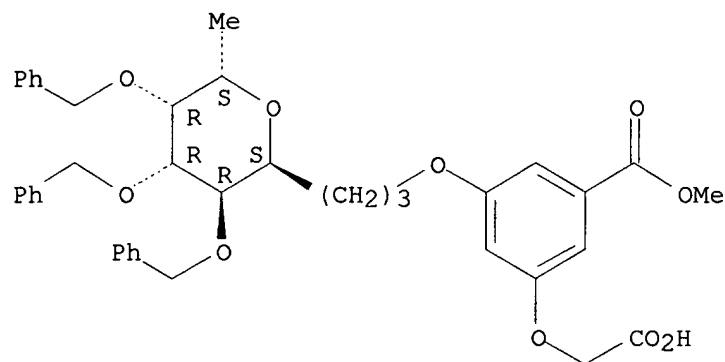
For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

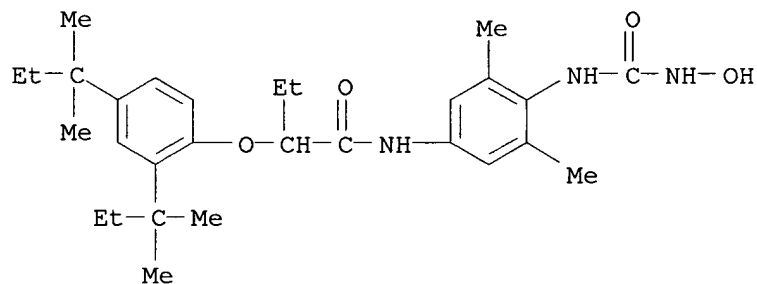
L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN L-glycero-D-galacto-Nonitol, 2,6-anhydro-9-O-[3-(carboxymethoxy)-5-(methoxycarbonyl)phenyl]-1,7,8-trideoxy-3,4,5-tris-O-(phenylmethyl)- (9CI)
MF C40 H44 O10

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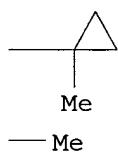
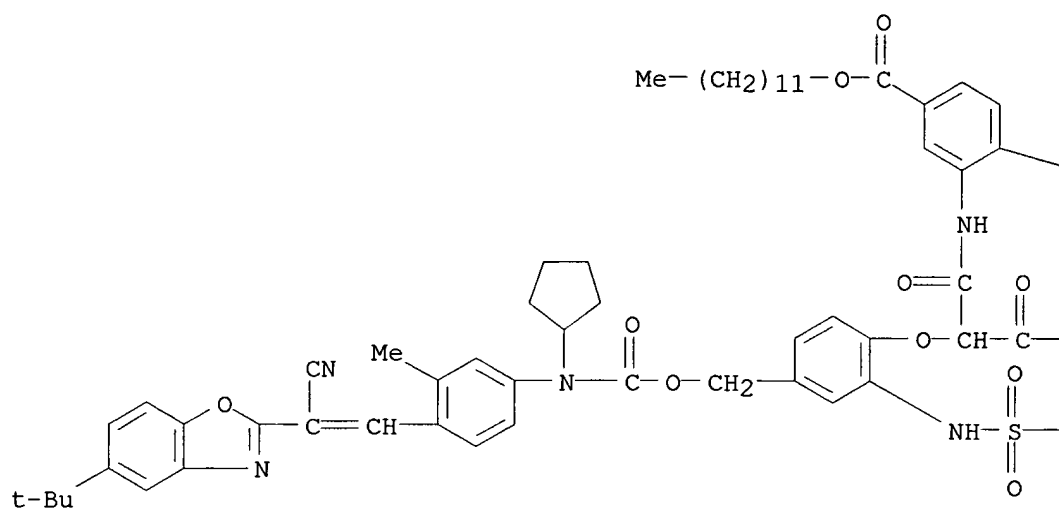
Absolute stereochemistry.



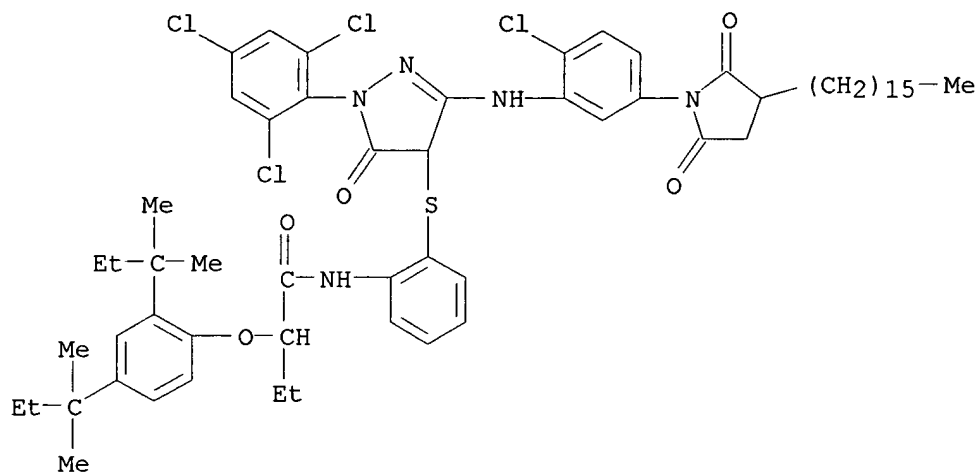
L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[4-
[[(hydroxyamino) carbonyl] amino]-3,5-dimethylphenyl]- (9CI)
MF C29 H43 N3 O4



L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Benzoic acid, 4-chloro-3-[[2-[4-[[[[4-[2-cyano-2-[5-(1,1-dimethylethyl)-2-benzoxazolyl]ethenyl]-3-methylphenyl]cyclopentylamino]carbonyl]oxy]methyl]-2-[(methylsulfonyl)amino]phenoxy]-3-(1-methylcyclopropyl)-1,3-dioxopropyl]amino]-, dodecyl ester (9CI)
MF C61 H74 Cl N5 O10 S



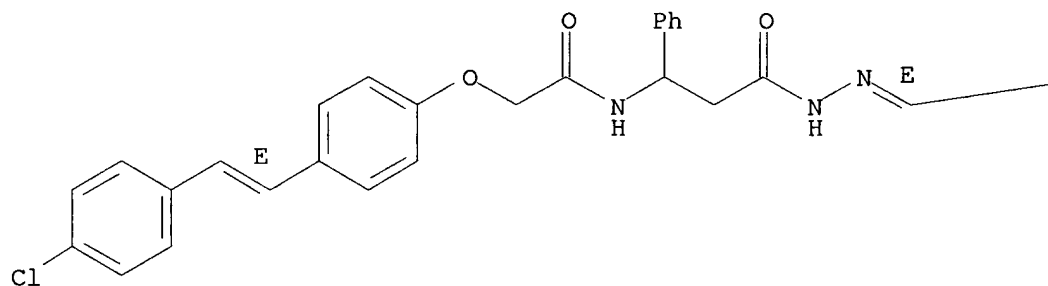
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L3 50 ANSWERS   REGISTRY   COPYRIGHT 2000 ACS
IN  Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[2-[[3-[[2-chloro-5-
    (3-hexadecyl-2,5-dioxo-1-pyrrolidinyl)phenyl]amino]-4,5-dihydro-5-oxo-1-
    (2,4,6-trichlorophenyl)-1H-pyrazol-4-yl]thio]phenyl]- (9CI)
MF  C61 H79 Cl4 N5 O5 S
```



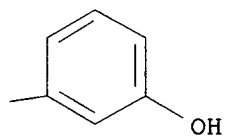
L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Benzenepropanoic acid, .beta.-[[[4-[(1E)-2-(4-chlorophenyl)ethenyl]phenoxy]acetyl]amino]-, (2E)-[(3-hydroxyphenyl)methylene]hydrazide (9CI)
 MF C32 H28 Cl N3 O4

Double bond geometry as shown.

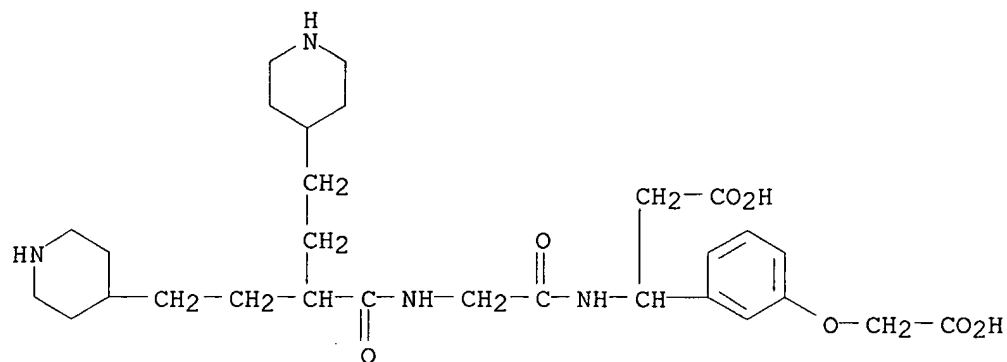
PAGE 1-A



PAGE 1-B

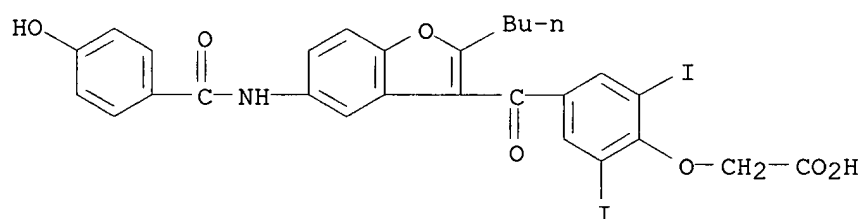


L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN .beta.-Alanine, N-[1-oxo-4-(4-piperidiny1)-2-[2-(4-piperidiny1)ethyl]butyl]glycyl-3-[3-(carboxymethoxy)phenyl]-, dihydrochloride (9CI)
 MF C29 H44 N4 O7 . 2 Cl H



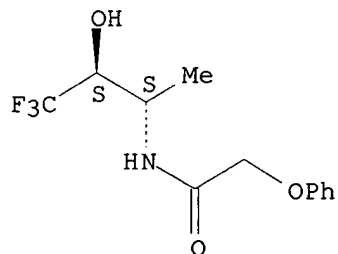
● 2 HCl

L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Acetic acid, [4-[[2-butyl-5-[(4-hydroxybenzoyl)amino]-3-benzofuranyl]carbonyl]-2,6-diiodophenoxy]- (9CI)
 MF C28 H23 I2 N O7



L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Acetamide, 2-phenoxy-N-(3,3,3-trifluoro-2-hydroxy-1-methylpropyl)-, [S-(R*,R*)]- (9CI)
 MF C12 H14 F3 N O3

Absolute stereochemistry.

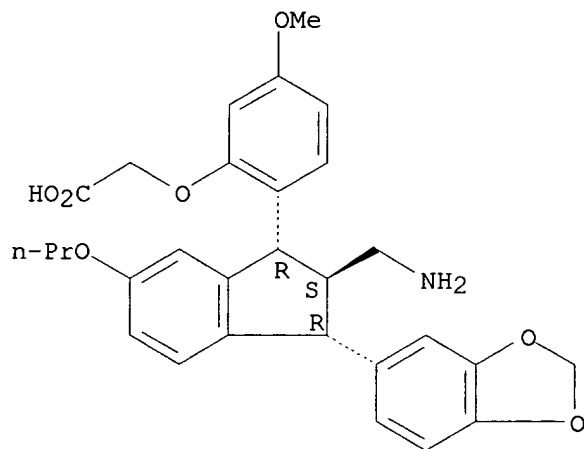


L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Acetic acid, [2-[(1R,2S,3R)-2-(aminomethyl)-3-(1,3-benzodioxol-5-yl)-2,3-

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MF C29 H31 N O7
dihydro-6-propoxy-1H-inden-1-yl]-5-methoxyphenoxy]-, rel- (9CI)

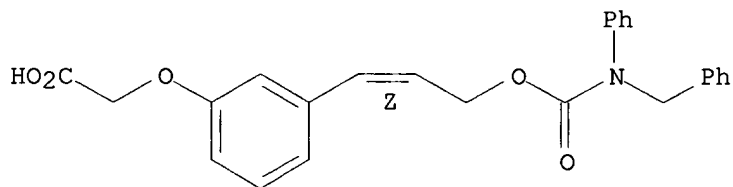
Relative stereochemistry.



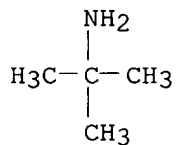
L3 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Acetic acid, [3-[(1Z)-3-[[[phenyl(phenylmethyl)amino]carbonyl]oxy]-1-propenyl]phenoxy]-, compd. with 2-methyl-2-propanamine (1:1) (9CI)
MF C25 H23 N O5 . C4 H11 N

CM 1

Double bond geometry as shown.



CM 2



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> log y

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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.90	1.05

STN INTERNATIONAL LOGOFF AT 15:33:56 ON 05 OCT 2000

Trying 3106016892...Open

Welcome to STN International! Enter x:x

LOGINID:sssptal611bxv

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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FILE 'HOME' ENTERED AT 15:36:51 ON 05 OCT 2000

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.15	0.15

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STRUCTURE FILE UPDATES: 4 OCT 2000 HIGHEST RN 292819-21-7
DICTIONARY FILE UPDATES: 4 OCT 2000 HIGHEST RN 292819-21-7

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 11, 2000

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT
for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 963

L1 SCREEN CREATED

=>

Uploading c:\stnexp4\queries\09543703.str

L2 STRUCTURE UPLOADED

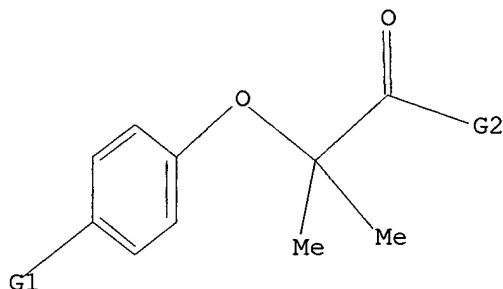
=> que L2 AND L1

L3 QUE L2 AND L1

=> d 12

L2 HAS NO ANSWERS

L2 STR



G1 C,H,O,N

G2 N,OH

Structure attributes must be viewed using STN Express query preparation.

=> s 12 sss sam

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SAMPLE SEARCH INITIATED 15:37:27 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 987 TO ITERATE

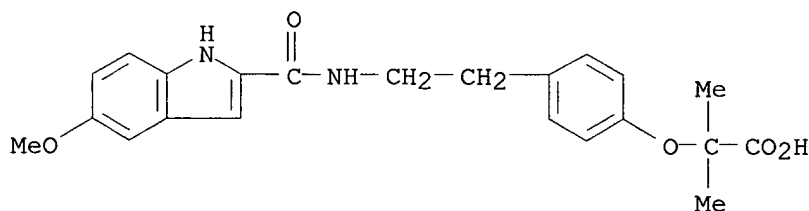
100.0% PROCESSED 987 ITERATIONS 50 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 17856 TO 21624
PROJECTED ANSWERS: 801 TO 1759

L4 50 SEA SSS SAM L2

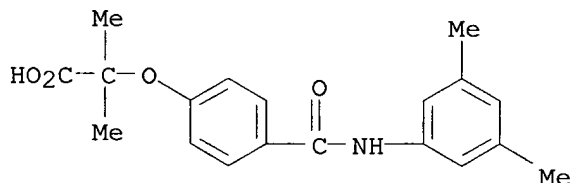
=> d scan 14

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[2-[[5-methoxy-1H-indol-2-yl)carbonyl]amino]ethyl]phenoxy]-2-methyl- (9CI)
MF C22 H24 N2 O5



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

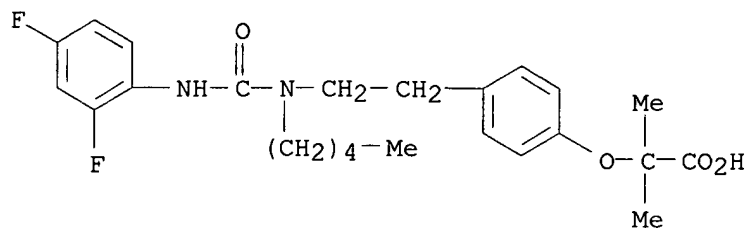
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[[3,5-dimethylphenyl)amino]carbonyl]phenoxy]-2-methyl- (9CI)
MF C19 H21 N O4



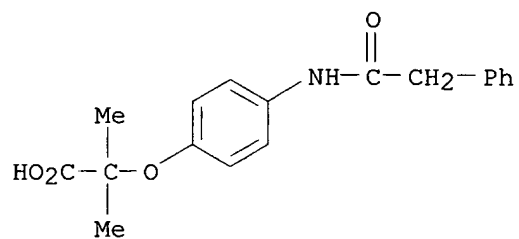
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[2-[[[2,4-difluorophenyl)amino]carbonyl]pentylamino]ethyl]phenoxy]-2-methyl- (9CI)
MF C24 H30 F2 N2 O4

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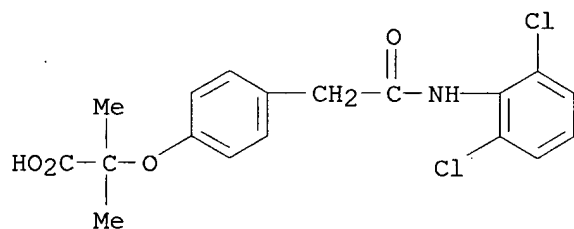


L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-methyl-2-[4-[(phenylacetyl)amino]phenoxy]-, monosodium salt (9CI)
MF C18 H19 N O4 . Na



● Na

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[2-[(2,6-dichlorophenyl)amino]-2-oxoethyl]phenoxy]-2-methyl-, monosodium salt (9CI)
MF C18 H17 Cl₂ N O₄ . Na



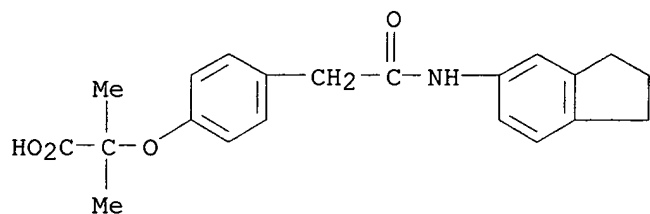
● Na

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[2-[(2,3-dihydro-1H-inden-5-yl)amino]-2-oxoethyl]phenoxy]-2-methyl- (9CI)

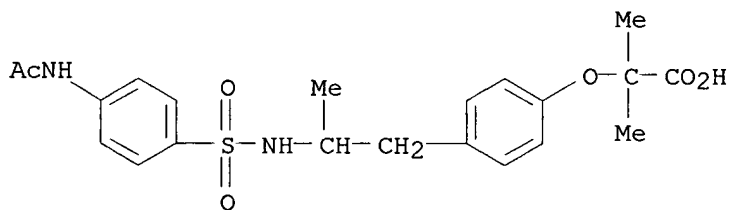
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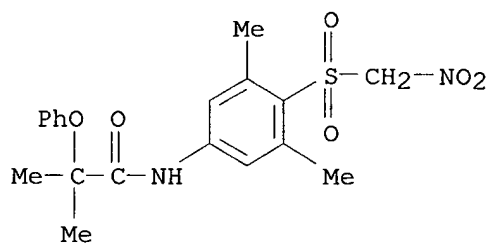
MF C21 H23 N O4
CI COM



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[2-[[[4-(acetylamino)phenyl]sulfonyl]amino]propyl]phenoxy]-2-methyl- (9CI)
MF C21 H26 N2 O6 S
CI COM

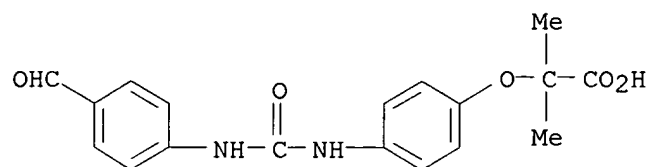


L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanamide, N-[3,5-dimethyl-4-[(nitromethyl)sulfonyl]phenyl]-2-methyl-2-phenoxy- (9CI)
MF C19 H22 N2 O6 S



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[[[(4-formylphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI)
MF C18 H18 N2 O5

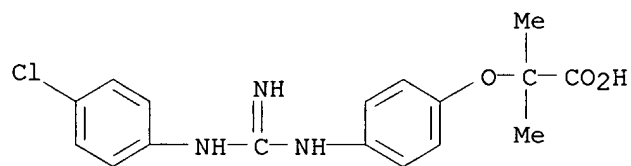
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L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN Propanoic acid, 2-[4-[[[(4-chlorophenyl)amino]iminomethyl]amino]phenoxy]-2-methyl- (9CI)

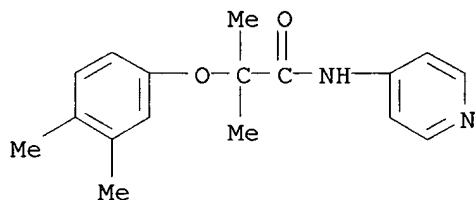
MF C17 H18 Cl N3 O3



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN Propanamide, 2-(3,4-dimethylphenoxy)-2-methyl-N-4-pyridinyl- (9CI)

MF C17 H20 N2 O2



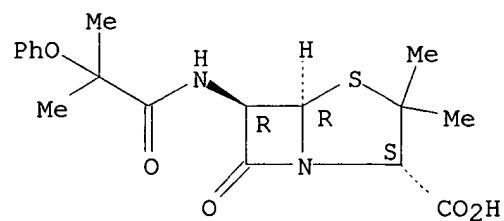
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 3,3-dimethyl-6-(2-methyl-2-phenoxypropionamido)-7-oxo-, sodium salt (7CI)

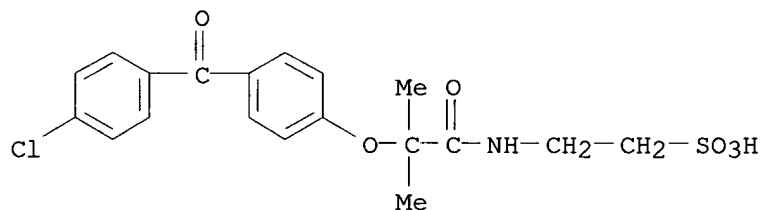
MF C18 H22 N2 O5 S . Na

Absolute stereochemistry.



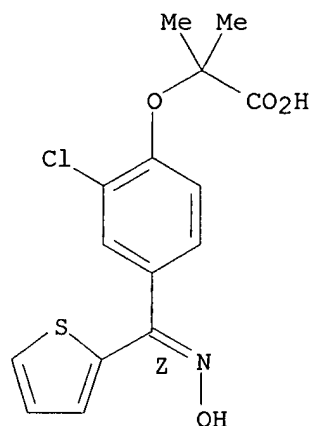
● Na

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Ethanesulfonic acid, 2-[[2-[4-(4-chlorobenzoyl)phenoxy]-2-methyl-1-oxopropyl]amino]- (9CI)
 MF C19 H20 Cl N O6 S
 CI COM

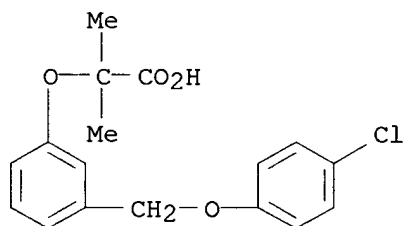


L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[2-chloro-4-[(hydroxyimino)-2-thienylmethyl]phenoxy]-2-methyl-, (Z)- (9CI)
 MF C15 H14 Cl N O4 S

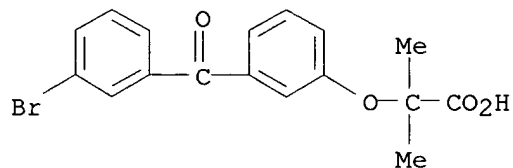
Double bond geometry as shown.



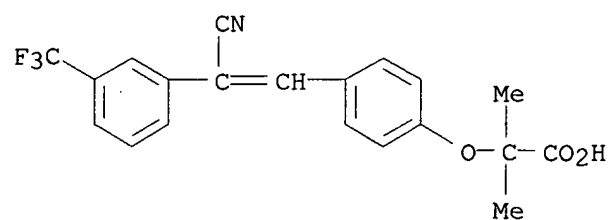
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[3-[(4-chlorophenoxy)methyl]phenoxy]-2-methyl- (9CI)
 MF C17 H17 Cl O4



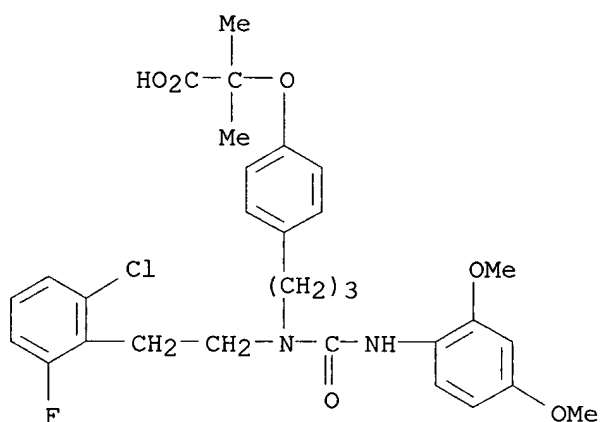
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[3-(3-bromobenzoyl)phenoxy]-2-methyl- (9CI)
 MF C17 H15 Br O4



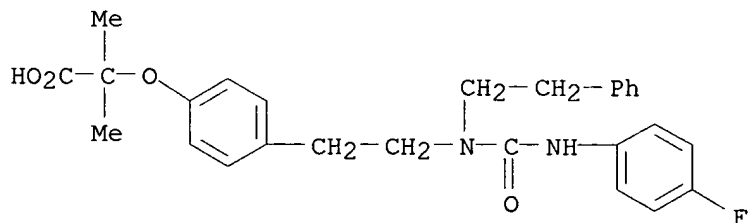
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[4-[2-cyano-2-[3-(trifluoromethyl)phenyl]ethenyl]phenoxy]-2-methyl- (9CI)
 MF C20 H16 F3 N O3



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[4-[3-[[2-(2-chloro-6-fluorophenyl)ethyl][[(2,4-dimethoxyphenyl)amino]carbonyl]amino]propyl]phenoxy]-2-methyl- (9CI)
 MF C30 H34 Cl F N2 O6



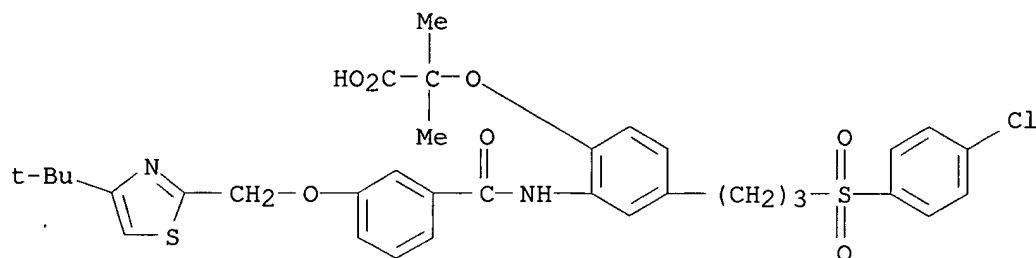
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[4-[2-[[[(4-fluorophenyl)amino]carbonyl](2-phenylethyl)amino]ethyl]phenoxy]-2-methyl- (9CI)
 MF C27 H29 F N2 O4



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[4-[3-[[4-(1,1-dimethylethyl)-2-thiazolyl]methoxy]benzoyl]amino]phenoxy]-2-methyl- (9CI)

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MF C34 H37 Cl N2 O7 S2

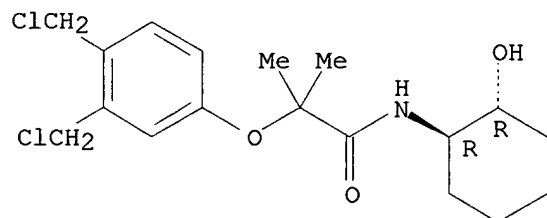


L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN Propanamide, 2-[3,4-bis(chloromethyl)phenoxy]-N-(2-hydroxycyclohexyl)-2-methyl-, trans- (9CI)

MF C18 H25 Cl2 N O3

Relative stereochemistry.

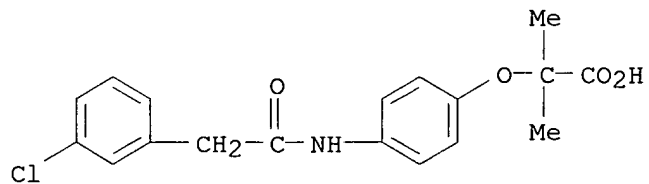


HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN Propanoic acid, 2-[4-[[3-chlorophenyl]acetyl]amino]phenoxy]-2-methyl-, monosodium salt (9CI)

MF C18 H18 Cl N O4 . Na



● Na

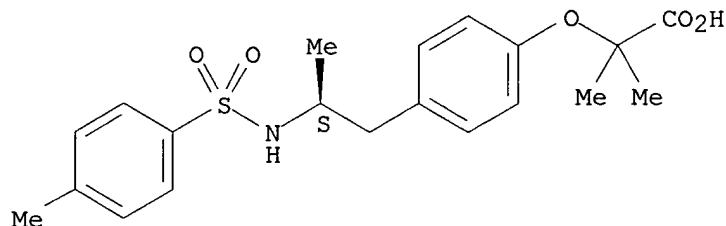
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS

09/825,925

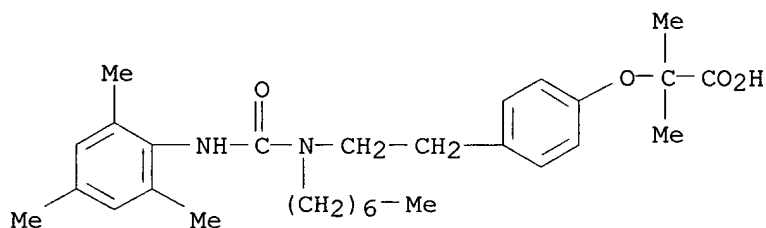
V. Balasubramanian

IN Propanoic acid, 2-methyl-2-[4-[2-[[4-methylphenyl)sulfonyl]amino]propyl]p
henoxy]-, (S)- (9CI)
MF C20 H25 N O5 S
CI COM

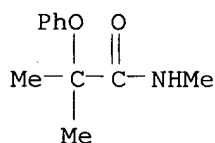
Absolute stereochemistry.



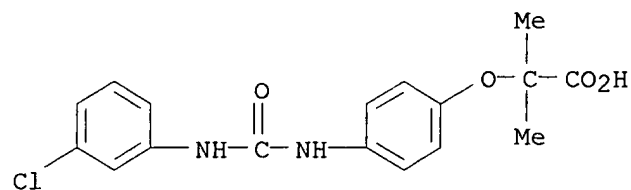
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[2-[heptyl[[2,4,6-trimethylphenyl)amino]carbonyl]ami
no]ethyl]phenoxy]-2-methyl- (9CI)
MF C29 H42 N2 O4



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanamide, N,2-dimethyl-2-phenoxy- (9CI)
MF C11 H15 N O2

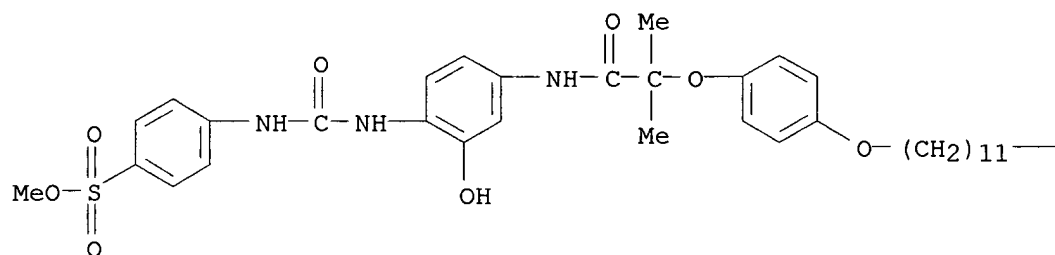


L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[[(3-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-
methyl- (9CI)
MF C17 H17 Cl N2 O4



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Benzenesulfonic acid, 4-[[[4-[[2-[4-(dodecyloxy)phenoxy]-2-methyl-1-oxopropyl]amino]-2-hydroxyphenyl]amino]carbonyl]amino]-, methyl ester (9CI)
 MF C36 H49 N3 O8 S

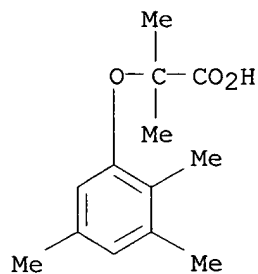
PAGE 1-A



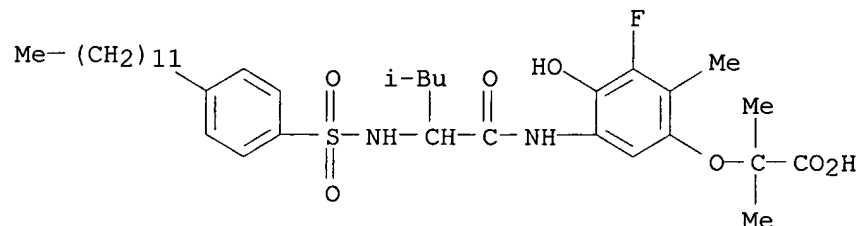
PAGE 1-B

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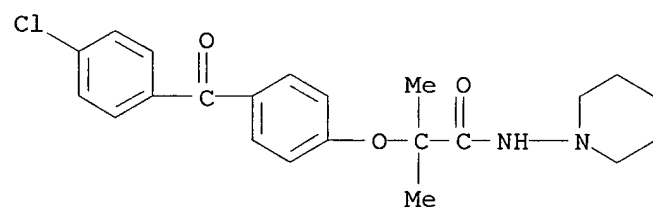
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-methyl-2-(2,3,5-trimethylphenoxy)- (9CI)
 MF C13 H18 O3



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[5-[[2-[[[(4-dodecylphenyl)sulfonyl]amino]-4-methyl-1-oxopentyl]amino]-3-fluoro-4-hydroxy-2-methylphenoxy]-2-methyl- (9CI)
 MF C35 H53 F N2 O7 S



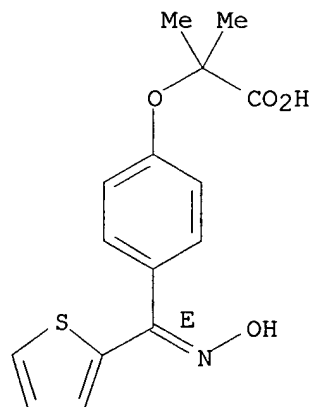
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanamide, 2-[4-(4-chlorobenzoyl)phenoxy]-2-methyl-N-1-piperidinyl- (9CI)
 MF C22 H25 Cl N2 O3



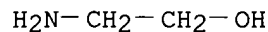
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-[4-[(hydroxyimino)-2-thienylmethyl]phenoxy]-2-methyl-, (E)-, compd. with 2-aminoethanol (1:1) (9CI)
 MF C15 H15 N O4 S . C2 H7 N O

CM 1

Double bond geometry as shown.

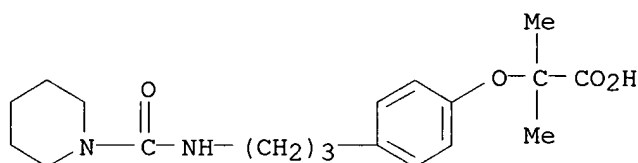


CM 2

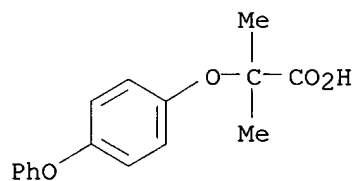


HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

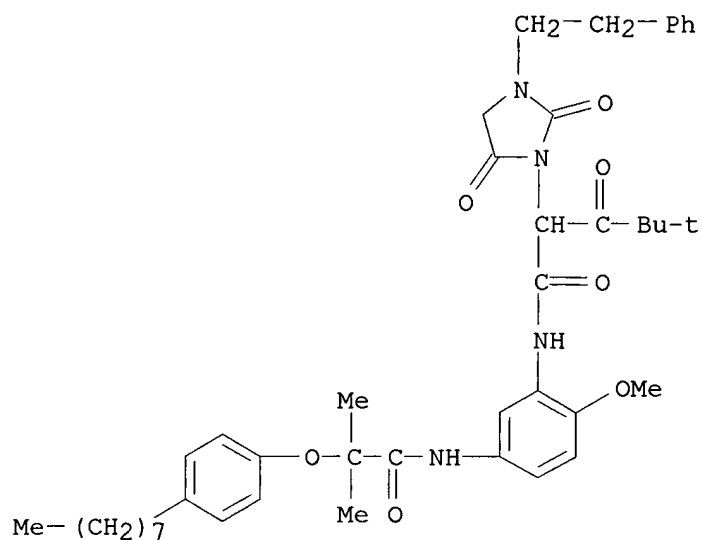
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-methyl-2-[4-[3-[(1-piperidinylcarbonyl)amino]propyl]phenoxy]- (9CI)
MF C19 H28 N2 O4
CI COM



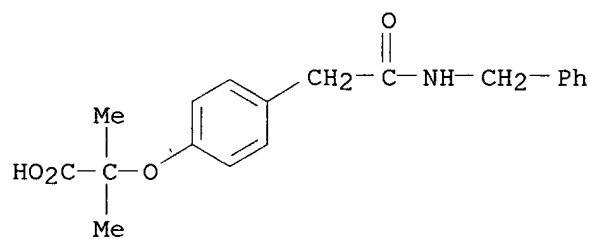
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-methyl-2-(4-phenoxyphenoxy)- (9CI)
MF C16 H16 O4



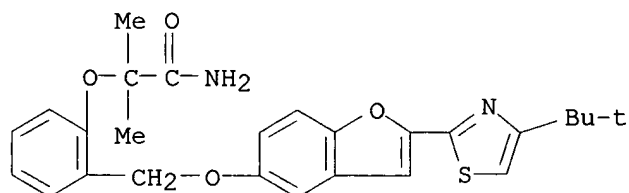
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN 1-Imidazolidineacetamide, .alpha.-(2,2-dimethyl-1-oxopropyl)-N-[2-methoxy-5-[[2-methyl-2-(4-octylphenoxy)-1-oxopropyl]amino]phenyl]-2,5-dioxo-3-(2-phenylethyl)- (9CI)
MF C43 H56 N4 O7



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanoic acid, 2-methyl-2-[4-[2-oxo-2-[(phenylmethyl)amino]ethyl]phenoxy]-
 (9CI)
 MF C19 H21 N O4



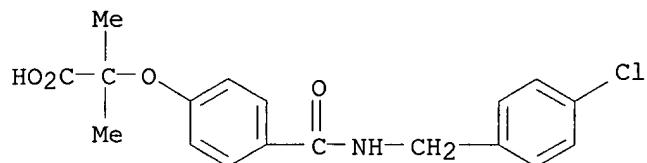
L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
 IN Propanamide, 2-[2-[[[2-[4-(1,1-dimethylethyl)-2-thiazolyl]-5-benzofuranyl]oxy]methyl]phenoxy]-2-methyl- (9CI)
 MF C26 H28 N2 O4 S



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS

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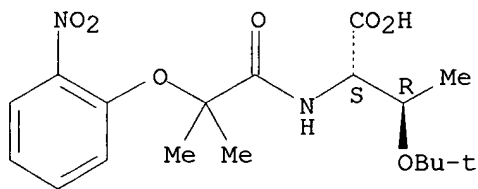
IN Propanoic acid, 2-[4-[[[(4-chlorophenyl)methyl]amino]carbonyl]phenoxy]-2-methyl-, monosodium salt (9CI)
MF C18 H18 Cl N O4 . Na



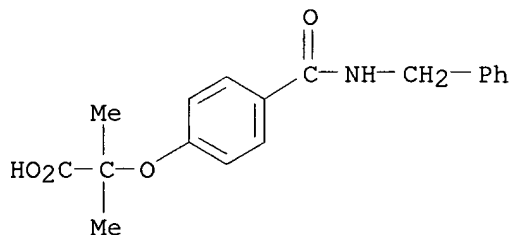
● Na

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN L-Threonine, O-(1,1-dimethylethyl)-N-[2-methyl-2-(2-nitrophenoxy)-1-oxopropyl]- (9CI)
MF C18 H26 N2 O7

Absolute stereochemistry.



L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-methyl-2-[4-[[[(phenylmethyl)amino]carbonyl]phenoxy]- (9CI)
MF C18 H19 N O4
CI COM

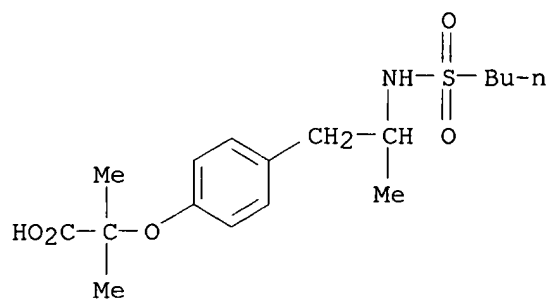


L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[2-[(butylsulfonyl)amino]propyl]phenoxy]-2-methyl-, monosodium salt (9CI)

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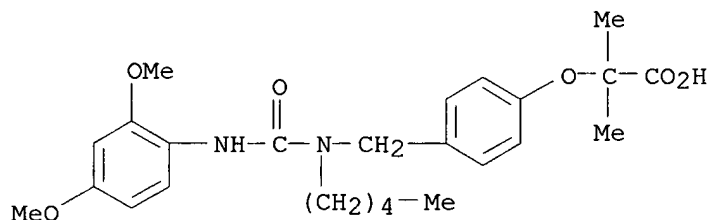
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MF C17 H27 N O5 S . Na



● Na

L4 50 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[[[[(2,4-dimethoxyphenyl)amino]carbonyl]pentylamino]methyl]phenoxy]-2-methyl- (9CI)
MF C25 H34 N2 O6



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.90

1.05

STN INTERNATIONAL LOGOFF AT 15:38:41 ON 05 OCT 2000

Trying 3106016892...Open

Welcome to STN International! Enter x:x

LOGINID:sssptal611bxv

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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DIOGENES
NEWS 3 Aug 21 CAS patent coverage expanded
NEWS 4 Aug 24 TABULATE Now Available in More STN Databases
NEWS 5 Aug 28 MEDLINE from 1958 to Date - Only on STN
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NEWS 7 Sep 11 Textile Technology Digest (TEXTILETECH) now available
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NEWS 9 Sep 29 The Philippines Inventory of Chemicals and Chemical
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FILE 'HOME' ENTERED AT 15:52:45 ON 05 OCT 2000

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.15	0.15

FILE 'REGISTRY' ENTERED AT 15:52:50 ON 05 OCT 2000

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STRUCTURE FILE UPDATES: 4 OCT 2000 HIGHEST RN 292819-21-7
DICTIONARY FILE UPDATES: 4 OCT 2000 HIGHEST RN 292819-21-7

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 11, 2000

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT

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for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 963

L1 SCREEN CREATED

=>

Uploading c:\stnexp4\queries\09543703.str

L2 STRUCTURE UPLOADED

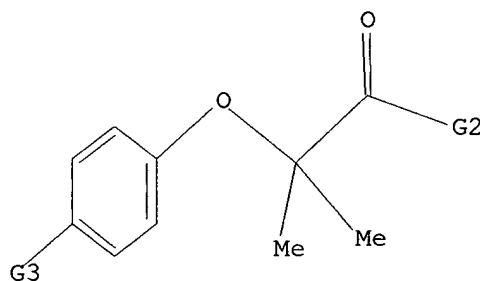
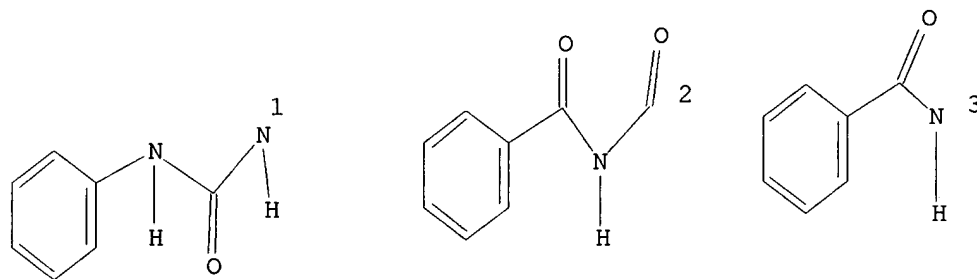
=> que L2 AND L1

L3 QUE L2 AND L1

=> d 12

L2 HAS NO ANSWERS

L2 STR



G1 C,H,O,N

G2 N,OH

G3 [01],[02],[03]

Structure attributes must be viewed using STN Express query preparation.

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=> s l2 sss sam

SAMPLE SEARCH INITIATED 15:53:35 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 7 TO ITERATE

100.0% PROCESSED 7 ITERATIONS
SEARCH TIME: 00.00.01

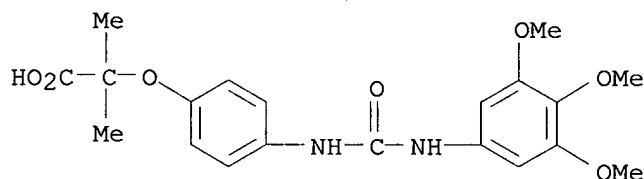
3 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 7 TO 298
PROJECTED ANSWERS: 3 TO 163

L4 3 SEA SSS SAM L2

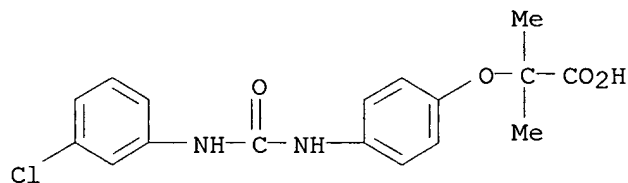
=> d scan l4

L4 3 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trimethoxyphenyl) amino] carbonyl] amino]phenoxy]- (9CI)
MF C20 H24 N2 O7



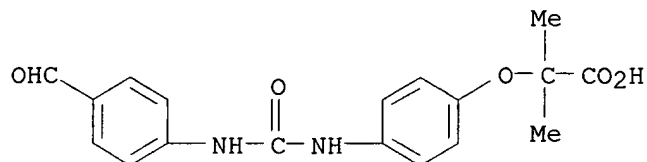
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L4 3 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[[[(3-chlorophenyl) amino] carbonyl] amino]phenoxy]-2-methyl- (9CI)
MF C17 H17 Cl N2 O4



L4 3 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Propanoic acid, 2-[4-[[[(4-formylphenyl) amino] carbonyl] amino]phenoxy]-2-methyl- (9CI)
MF C18 H18 N2 O5

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ALL ANSWERS HAVE BEEN SCANNED

=> s l2 sss ful

FULL SEARCH INITIATED 15:54:07 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 136 TO ITERATE

100.0% PROCESSED 136 ITERATIONS 63 ANSWERS
SEARCH TIME: 00.00.01

L5 63 SEA SSS FUL L2

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	126.60	126.75

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FILE COVERS 1967 - 5 Oct 2000 VOL 133 ISS 15
FILE LAST UPDATED: 4 Oct 2000 (20001004/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

This file supports REGISTRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

Now you can extend your author, patent assignee, patent information, and title searches back to 1907. The records from 1907-1966 now have this searchable data in CAOLD. You now have electronic access to all of CA: 1907 to 1966 in CAOLD and 1967 to the present in CAPLUS on STN.

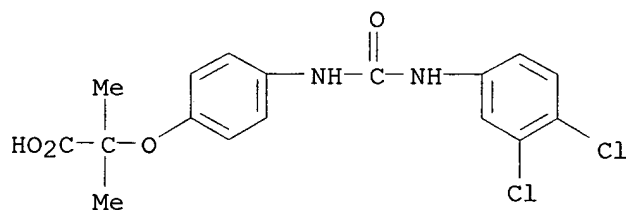
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L6 19 L5

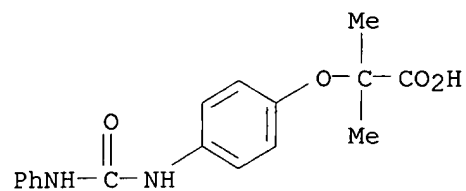
09/825,925

=> d l6 1-19 bib hitstr

L6 ANSWER 1 OF 19 CAPLUS COPYRIGHT 2000 ACS
AN 1999:697456 CAPLUS
DN 132:44918
TI Novel Inhibitors of Advanced Glycation Endproducts. [Erratum to document cited in CA131:317717]
AU Rahbar, Samuel; Yernini, Kiran Kumar V.; Scott, Stephen; Gonzales, Noe; Lalezari, Iraj
CS Dep. Diabetes, Endocrinology & Metabolism, City of Hope National Medical Center, Duarte, CA, 91010-0269, USA
SO Biochem. Biophys. Res. Commun. (1999), 264(3), 1008
CODEN: BBRC99; ISSN: 0006-291X
PB Academic Press
DT Journal
LA English
IT 117011-50-4 121809-57-2 121809-58-3
121809-65-2 121809-67-4 121809-68-5
121809-69-6 121809-70-9 121809-73-2
121809-74-3 121809-75-4 121809-77-6
121809-80-1 121809-82-3 121809-83-4
121809-84-5 121809-85-6 121809-89-0
121809-90-3 121809-94-7 157580-09-1
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249513-69-7 249513-70-0 249513-71-1
249513-72-2 249513-73-3 249513-74-4
249513-76-6 249513-77-7 249513-80-2
249513-81-3 249513-83-5 249513-84-6
RL: BAC (Biological activity or effector, except adverse); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(novel inhibitors of advanced glycation endproducts (Erratum))
RN 117011-50-4 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

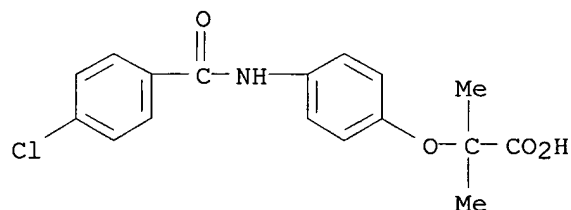


RN 121809-57-2 CAPLUS
CN Propanoic acid, 2-methyl-2-[4-[[[(phenylamino)carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



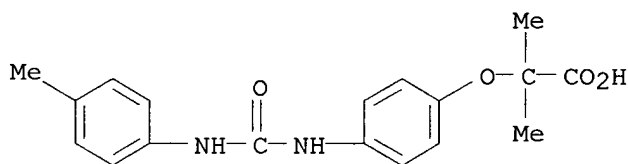
RN 121809-58-3 CAPLUS

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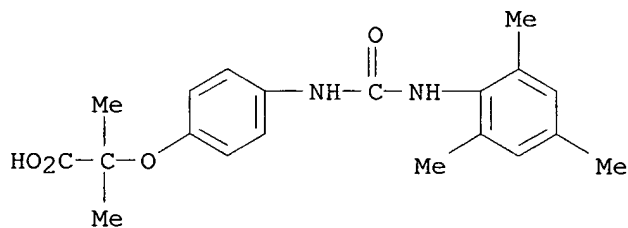
RN 121809-65-2 CAPLUS

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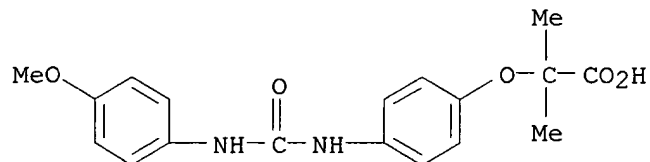
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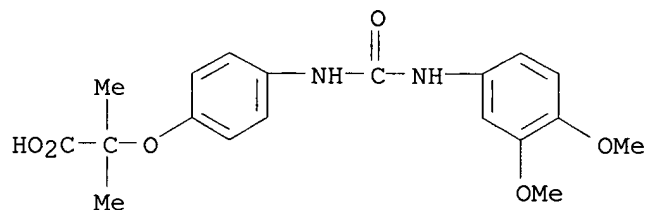
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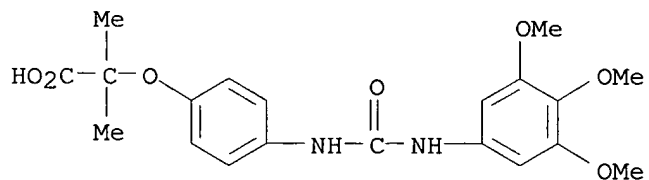
RN 121809-69-6 CAPLUS

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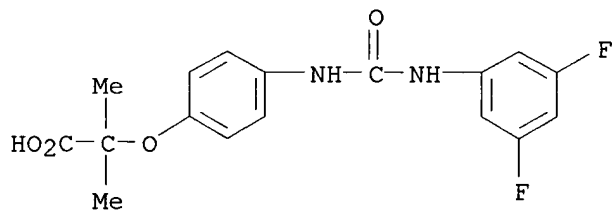
RN 121809-70-9 CAPLUS

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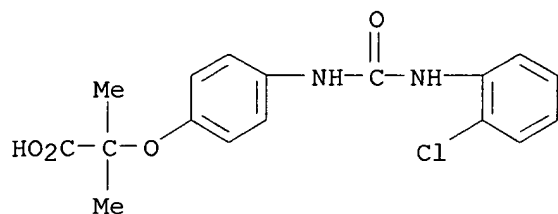
RN 121809-73-2 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-difluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



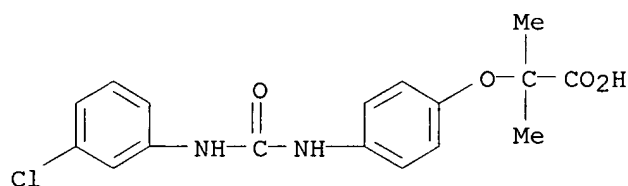
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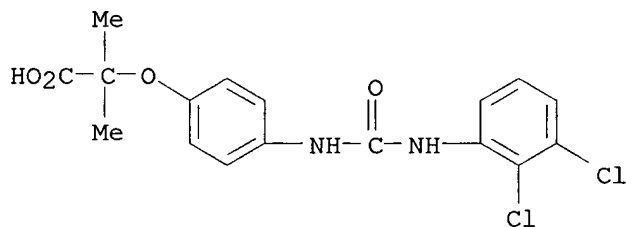
RN 121809-75-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



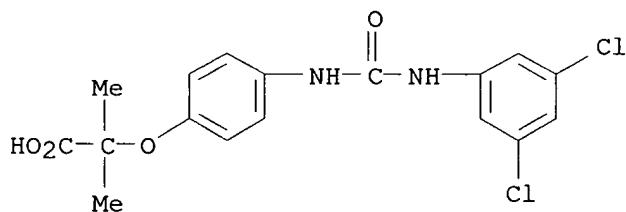
RN 121809-77-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(2,3-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



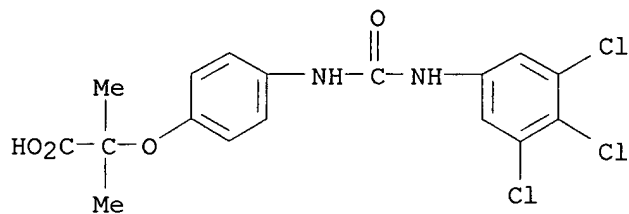
RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



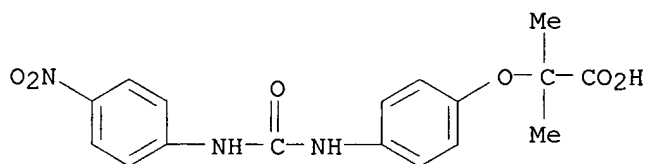
RN 121809-82-3 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



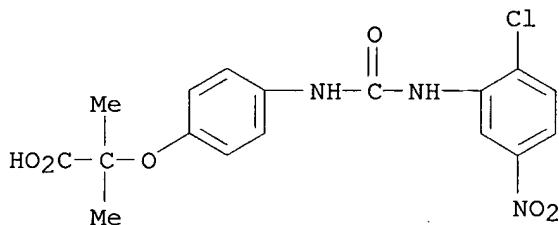
RN 121809-83-4 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(4-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



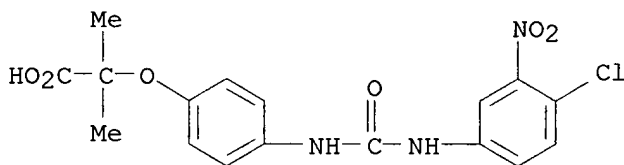
RN 121809-84-5 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-chloro-5-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



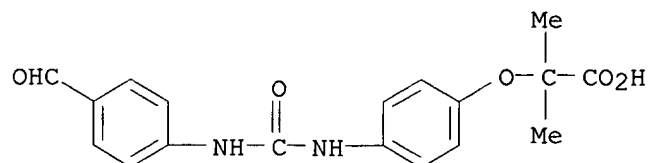
RN 121809-85-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-chloro-3-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



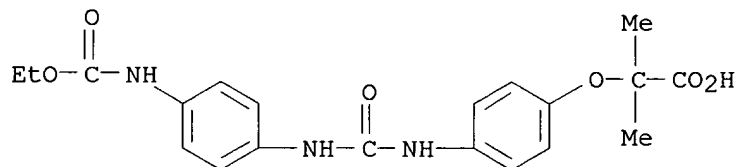
RN 121809-89-0 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-formylphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



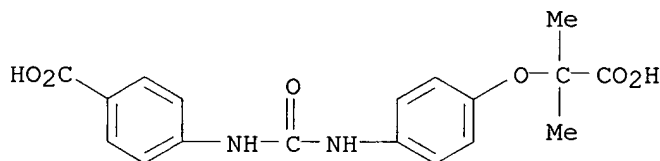
RN 121809-90-3 CAPLUS

CN Propanoic acid, 2-[4-[[[4-(ethoxycarbonyl)amino]phenyl]amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



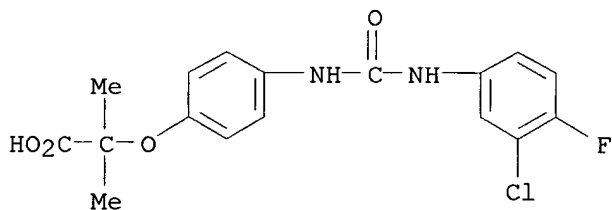
RN 121809-94-7 CAPLUS

CN Benzoic acid, 4-[[[4-(1-carboxy-1-methylethoxy)phenyl]amino]carbonyl]amino]phenyl]- (9CI) (CA INDEX NAME)



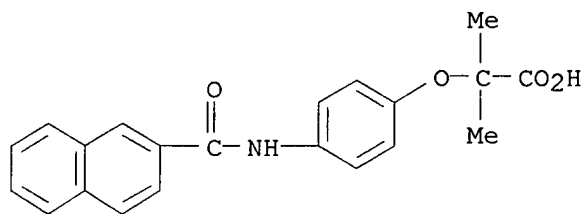
RN 157580-09-1 CAPLUS

CN Propanoic acid, 2-[4-[[[3-chloro-4-fluorophenyl]amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 249513-62-0 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[(2-naphthalenylcarbonyl)amino]phenoxy]- (9CI) (CA INDEX NAME)

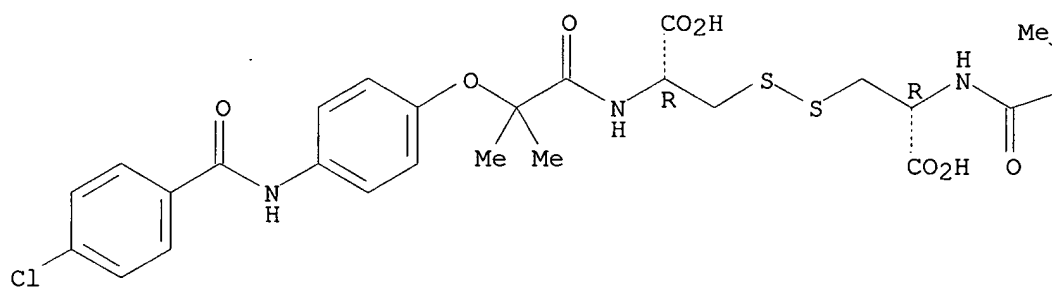


RN 249513-66-4 CAPLUS

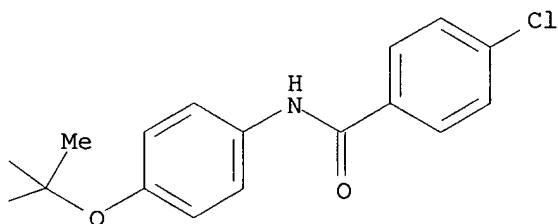
CN L-Cystine, N,N'-bis[2-[4-[(4-chlorobenzoyl)amino]phenoxy]-2-methyl-1-oxopropyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-A

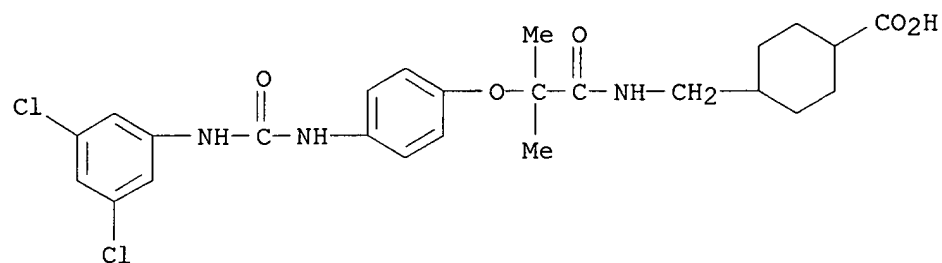


PAGE 1-B



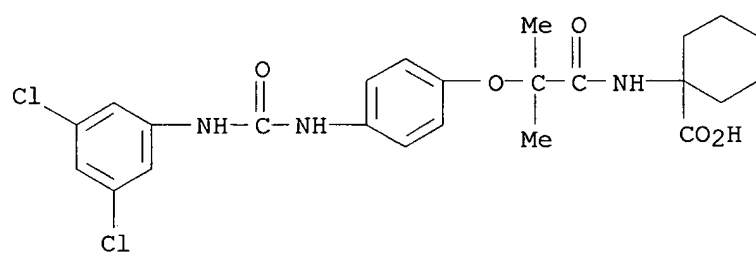
RN 249513-67-5 CAPLUS

CN Cyclohexanecarboxylic acid, 4-[[[2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl-1-oxopropyl]amino]methyl]- (9CI) (CA INDEX NAME)



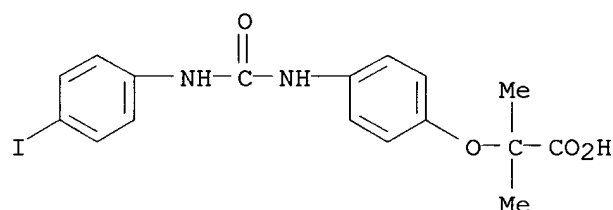
RN 249513-69-7 CAPLUS

CN Cyclohexanecarboxylic acid, 1-[[2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl-1-oxopropyl]amino]- (9CI) (CA INDEX NAME)



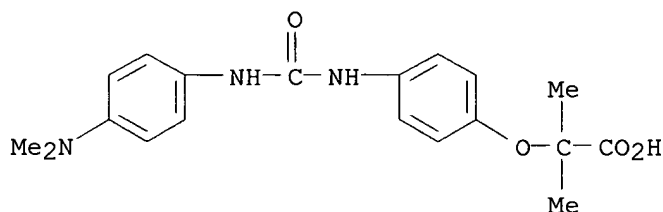
RN 249513-70-0 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-iodophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 249513-71-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-(dimethylamino)phenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

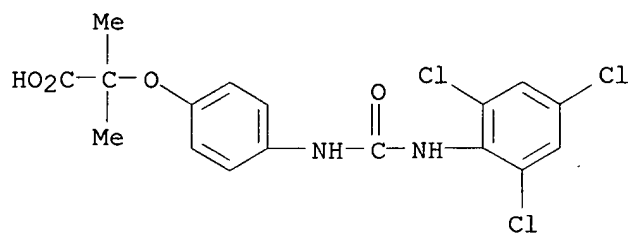


RN 249513-72-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,4,6-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)

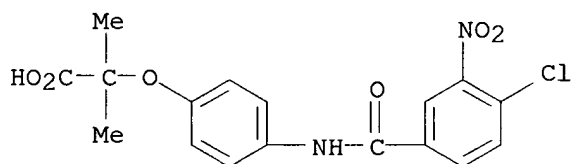
V. Balasubramanian

no]phenoxy]- (9CI) (CA INDEX NAME)



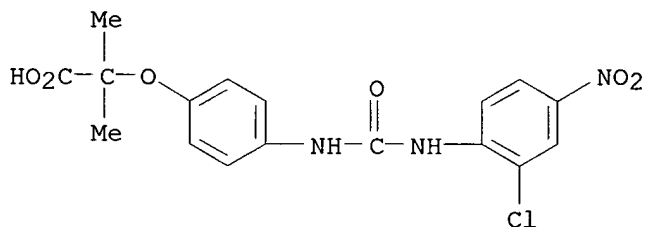
RN 249513-73-3 CAPLUS

CN Propanoic acid, 2-[4-[(4-chloro-3-nitrobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



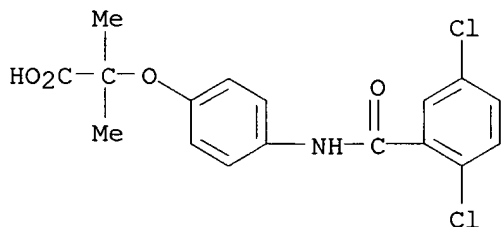
RN 249513-74-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-chloro-4-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 249513-76-6 CAPLUS

CN Propanoic acid, 2-[4-[(2,5-dichlorobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

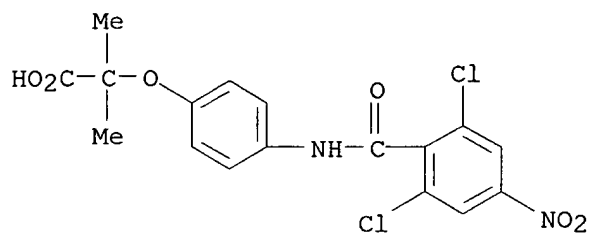


RN 249513-77-7 CAPLUS

CN Propanoic acid, 2-[4-[(2,6-dichloro-4-nitrobenzoyl)amino]phenoxy]-2-methyl-

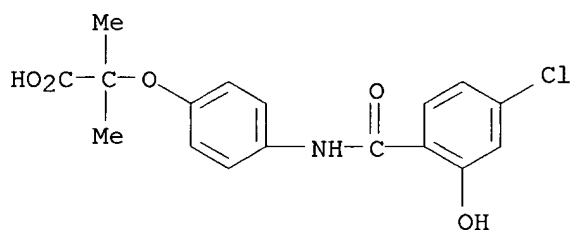
V. Balasubramanian

(9CI) (CA INDEX NAME)



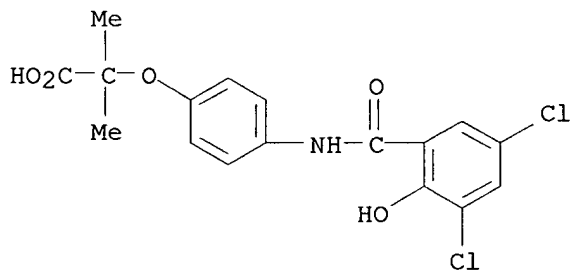
RN 249513-80-2 CAPLUS

CN Propanoic acid, 2-[4-[(4-chloro-2-hydroxybenzoyl)amino]phenoxy]-2-methyl-
(9CI) (CA INDEX NAME)



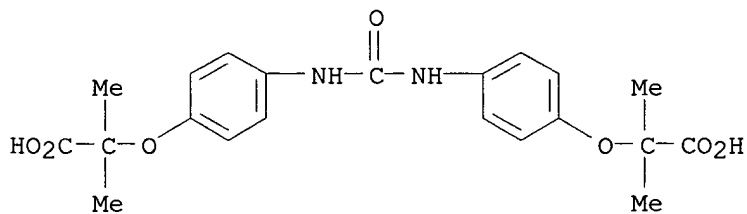
RN 249513-81-3 CAPLUS

CN Propanoic acid, 2-[4-[(3,5-dichloro-2-hydroxybenzoyl)amino]phenoxy]-2-methyl-
(9CI) (CA INDEX NAME)



RN 249513-83-5 CAPLUS

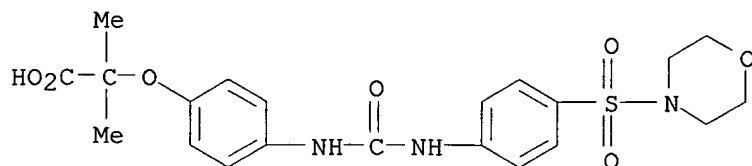
CN Propanoic acid, 2,2'-[carbonylbis(imino-4,1-phenyleneoxy)]bis[2-methyl-
(9CI) (CA INDEX NAME)



V. Balasubramanian

RN 249513-84-6 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[4-(4-morpholinylsulfonyl)phenyl]amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



L6 ANSWER 2 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1999:691824 CAPLUS

DN 132:81

TI A new rapid method to detect inhibition of Amadori product generated by .delta.-gluconolactone

AU Rahbar, Samuel; Nadler, Jerry L.

CS Department of Diabetes, Endocrinology and Metabolism, Gonda Diabetes Center, City of Hope National Medical Center, Duarte, CA, 91010, USA

SO Clin. Chim. Acta (1999), 287(1-2), 123-130

CODEN: CCATAR; ISSN: 0009-8981

PB Elsevier Science Ireland Ltd.

DT Journal

LA English

IT 121809-57-2 157580-09-1 249513-72-2

249513-74-4

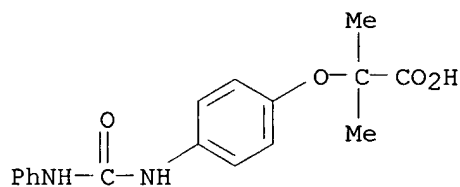
RL: BAC (Biological activity or effector, except adverse); THU

(Therapeutic use); BIOL (Biological study); USES (Uses)

(new rapid method to detect inhibition of Amadori product generated by .delta.-gluconolactone) 8

RN 121809-57-2 CAPLUS

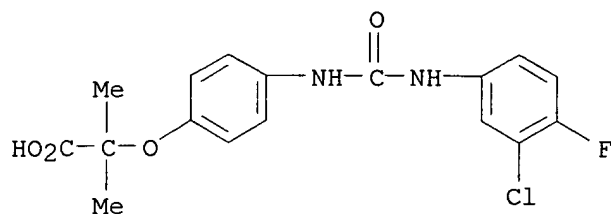
CN Propanoic acid, 2-methyl-2-[4-[(phenylamino)carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



RN 157580-09-1 CAPLUS

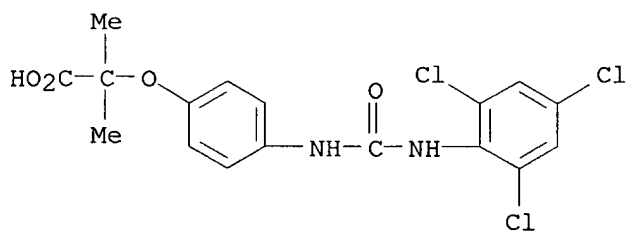
CN Propanoic acid, 2-[4-[[[(3-chloro-4-fluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

V. Balasubramanian



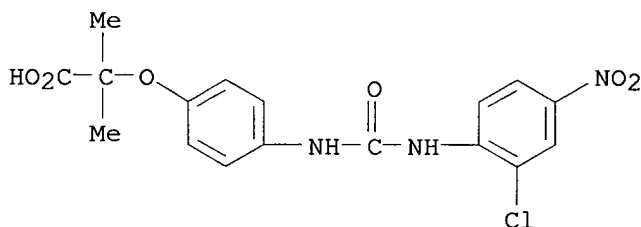
RN 249513-72-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,4,6-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



RN 249513-74-4 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2-chloro-4-nitrophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



RE.CNT 24

RE

- (1) Bailey, A; Mech Aging Dev 1998, V106, P1 CAPLUS
- (2) Baynes, J; Diabetes 1999, V48, P1 CAPLUS
- (4) Booth, A; J Biol Chem 1997, V272, P5430 CAPLUS
- (5) Brett, J; Am J Pathol 1993, V143, P1699 CAPLUS
- (6) Bucala, R; Endocrinology cardiovascular function 1998, P159 CAPLUS

ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 3 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1999:635489 CAPLUS

DN 131:262584

TI Modified hemoglobin and its use as a component of an artificial blood substitute

IN Lalezari, Iraj; Lalezari, Parviz

PA Montefiore Medical Center, USA

SO U.S., 22 pp., Cont. of U.S. Ser. No. 380,097, abandoned.

CODEN: USXXAM

09/825,925

V. Balasubramanian

DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 5962651	A	19991005	US 1997-903930	19970731
	US 6072072	A	20000606	US 1999-274072	19990322

PRAI US 1995-380097 19950127
US 1997-903930 19970731

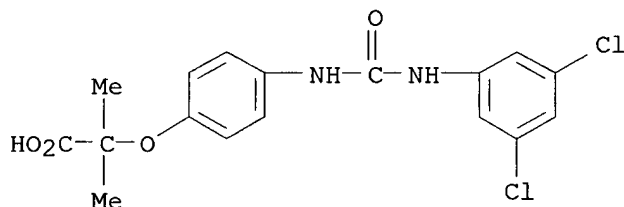
OS MARPAT 131:262584

IT **121809-80-1P 245075-84-7P 245075-86-9P**
245075-87-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(prepn. of functional compds. for Hb modification for blood
substitutes)

RN 121809-80-1 CAPLUS

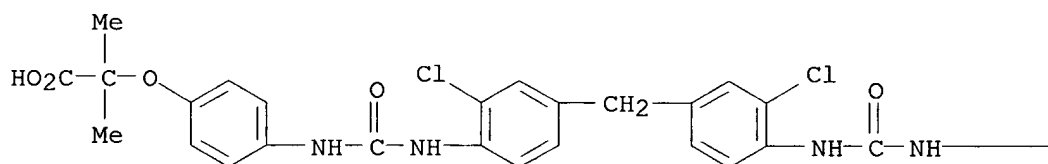
CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



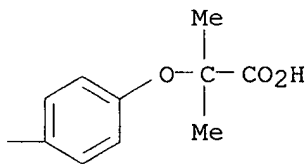
RN 245075-84-7 CAPLUS

CN Propanoic acid, 2,2'-[methylenebis[(2-chloro-4,1-
phenylene)iminocarbonylimino-4,1-phenyleneoxy]]bis[2-methyl- (9CI) (CA
INDEX NAME)

PAGE 1-A



PAGE 1-B



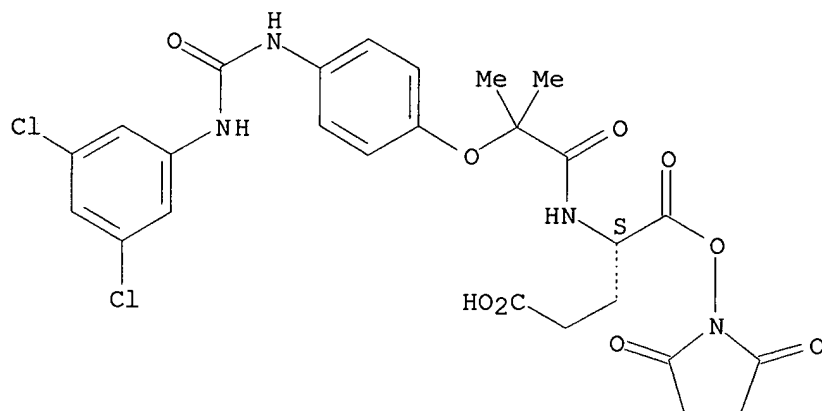
RN 245075-86-9 CAPLUS

CN Pentanoic acid, 4-[[2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl-1-oxopropyl]amino]-5-[(2,5-dioxo-1-pyrrolidinyl)oxy]-5-oxo-,
(4S)- (9CI) (CA INDEX NAME)

09/825,925

V. Balasubramanian

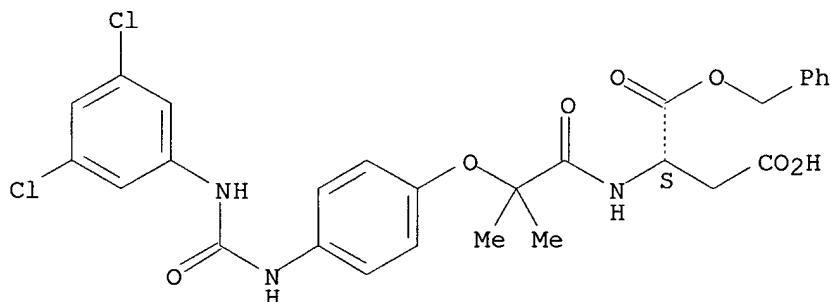
Absolute stereochemistry.



RN 245075-87-0 CAPLUS

CN L-Aspartic acid, N-[2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl-1-oxopropyl]-, 1-(phenylmethyl) ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RE.CNT 4

RE

(1) Abraham; US 5122539 1992 CAPLUS

(2) Calezari; US 5268500 1993

(3) Fronticelli; US 5239061 1993

(4) Hoffman; US 5028588 1991

L6 ANSWER 4 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1999:558280 CAPLUS

DN 131:317717

TI Novel Inhibitors of Advanced Glycation Endproducts

AU Rahbar, Samuel; Kumar Yernini, Kiran; Scott, Stephen; Gonzales, Noe; Lalezari, Iraj

CS Department of Diabetes, Endocrinology & Metabolism, City of Hope National Medical Center, Duarte, CA, 91010-0269, USA

SO Biochem. Biophys. Res. Commun. (1999), 262(3), 651-656

CODEN: BBRC9; ISSN: 0006-291X

PB Academic Press

DT Journal

LA English

09/825,925

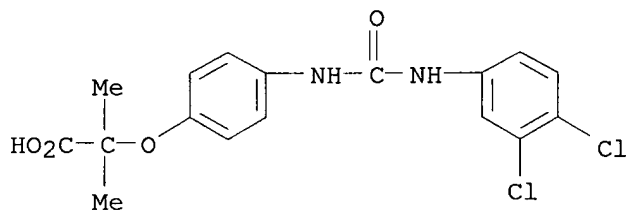
V. Balasubramanian

IT 117011-50-4 121809-57-2 121809-58-3
121809-65-2 121809-67-4 121809-68-5
121809-69-6 121809-70-9 121809-73-2
121809-74-3 121809-75-4 121809-77-6
121809-80-1 121809-82-3 121809-83-4
121809-84-5 121809-85-6 121809-89-0
121809-90-3 121809-94-7 157580-09-1
249513-62-0 249513-66-4 249513-67-5
249513-69-7 249513-70-0 249513-71-1
249513-72-2 249513-73-3 249513-74-4
249513-76-6 249513-77-7 249513-80-2
249513-81-3 249513-83-5 249513-84-6

RL: BAC (Biological activity or effector, except adverse); THU
(Therapeutic use); BIOL (Biological study); USES (Uses)
(novel inhibitors of advanced glycation endproducts)

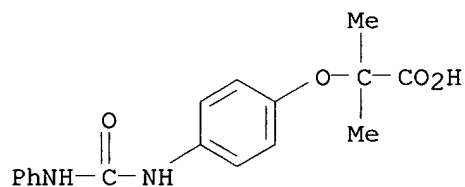
RN 117011-50-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



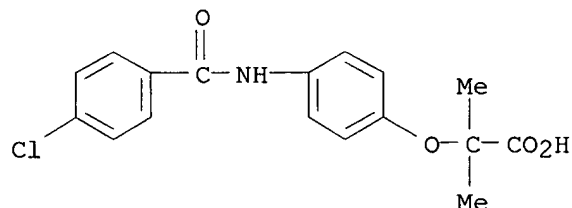
RN 121809-57-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[(phenylamino) carbonyl] amino]phenoxy]-
(9CI) (CA INDEX NAME)



RN 121809-58-3 CAPLUS

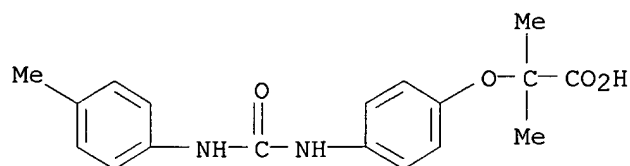
CN Propanoic acid, 2-[4-[(4-chlorobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA
INDEX NAME)



RN 121809-65-2 CAPLUS

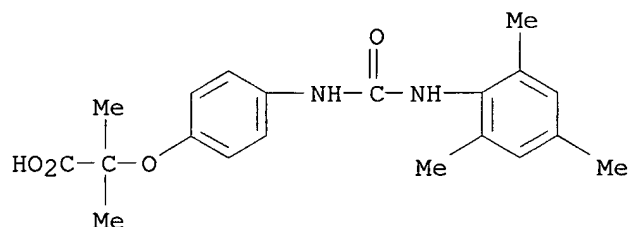
CN Propanoic acid, 2-methyl-2-[4-[[[(4-methylphenyl)amino]carbonyl]amino]phen

oxy]- (9CI) (CA INDEX NAME)



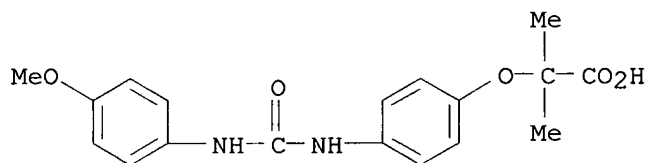
RN 121809-67-4 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,4,6-trimethylphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



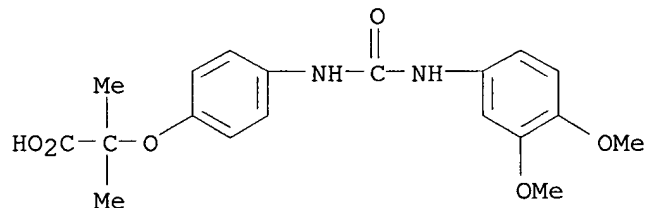
RN 121809-68-5 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-methoxyphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-69-6 CAPLUS

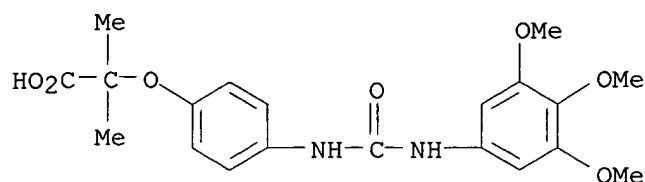
CN Propanoic acid, 2-[4-[[[(3,4-dimethoxyphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-70-9 CAPLUS

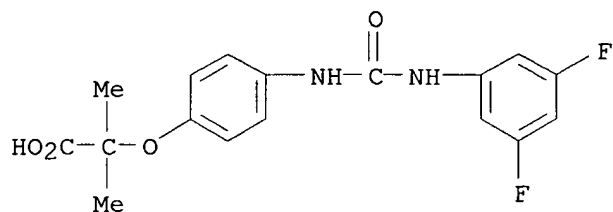
CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trimethoxyphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

V. Balasubramanian



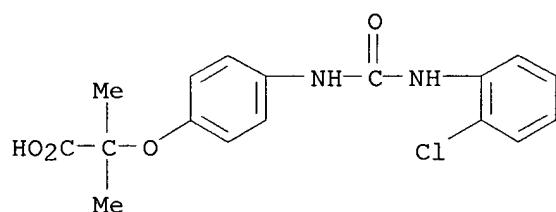
RN 121809-73-2 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-difluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



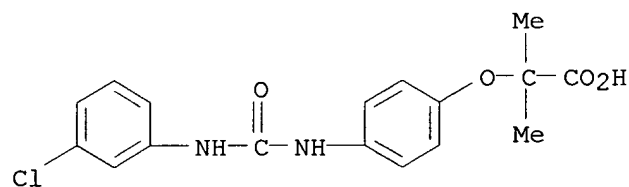
RN 121809-74-3 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



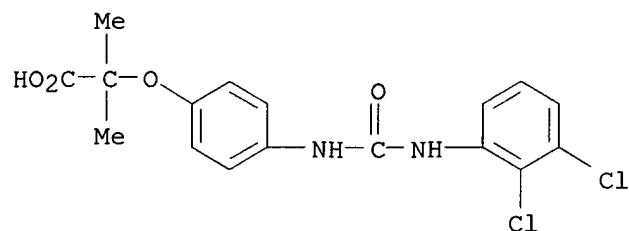
RN 121809-75-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



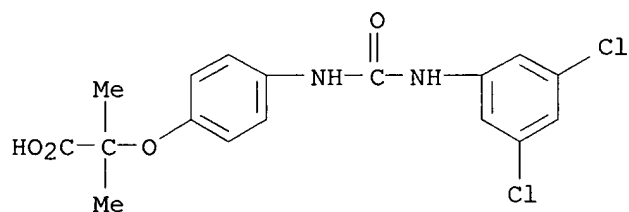
RN 121809-77-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(2,3-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



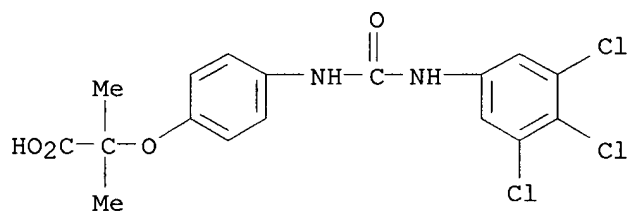
RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



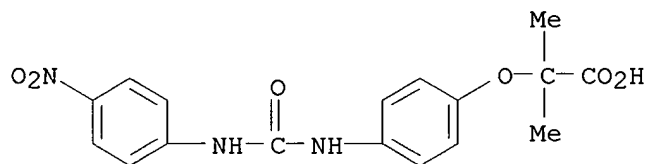
RN 121809-82-3 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



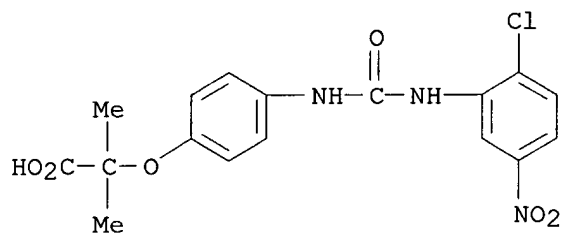
RN 121809-83-4 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(4-nitrophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



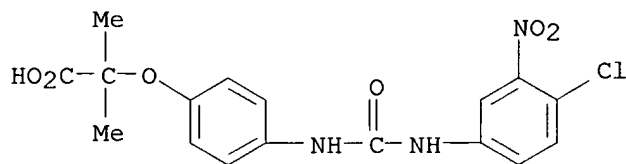
RN 121809-84-5 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-chloro-5-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



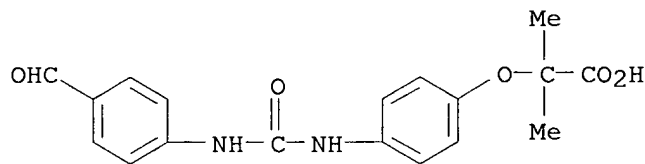
RN 121809-85-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-chloro-3-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



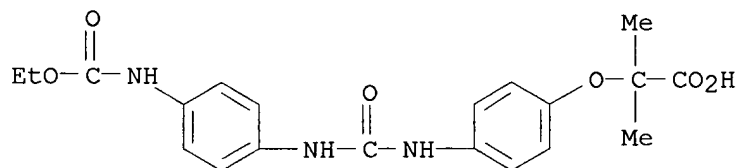
RN 121809-89-0 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-formylphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-90-3 CAPLUS

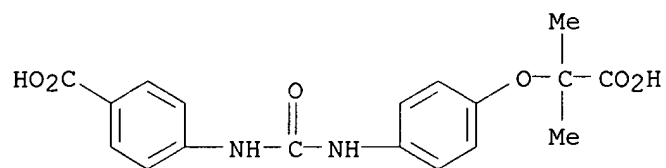
CN Propanoic acid, 2-[4-[[[4-[(ethoxycarbonyl)amino]phenyl]amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-94-7 CAPLUS

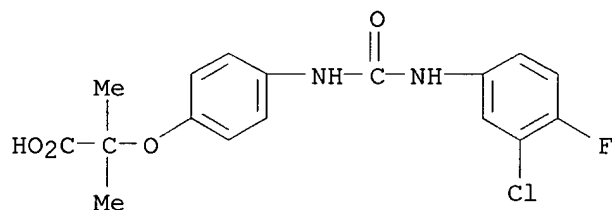
CN Benzoic acid, 4-[[[4-[(1-carboxy-1-methylethoxy)phenyl]amino]carbonyl]amino]phenyl- (9CI) (CA INDEX NAME)

V. Balasubramanian



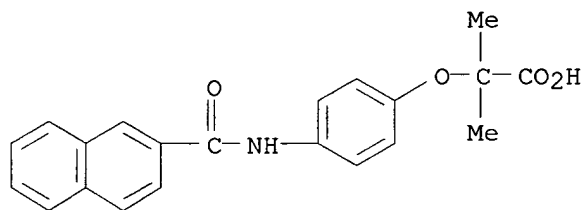
RN 157580-09-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-chloro-4-fluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 249513-62-0 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[(2-naphthalenylcarbonyl)amino]phenoxy]- (9CI) (CA INDEX NAME)

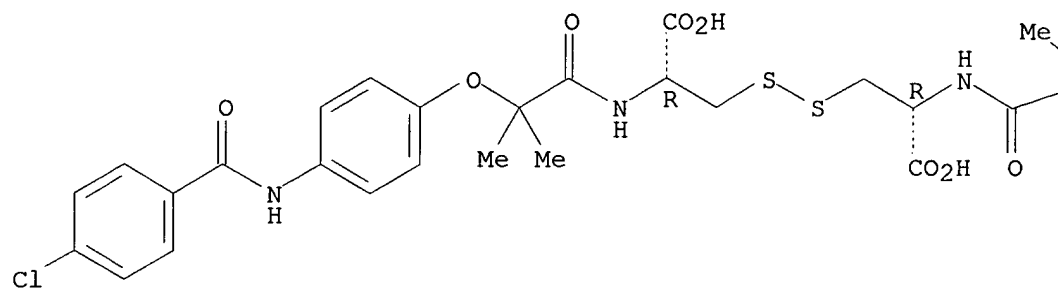


RN 249513-66-4 CAPLUS

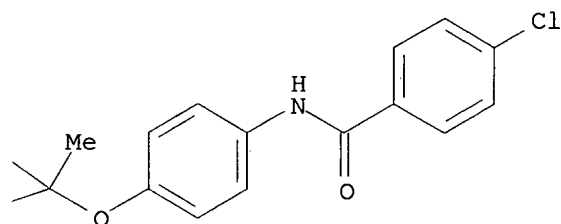
CN L-Cystine, N,N'-bis[2-[4-[(4-chlorobenzoyl)amino]phenoxy]-2-methyl-1-oxopropyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-A

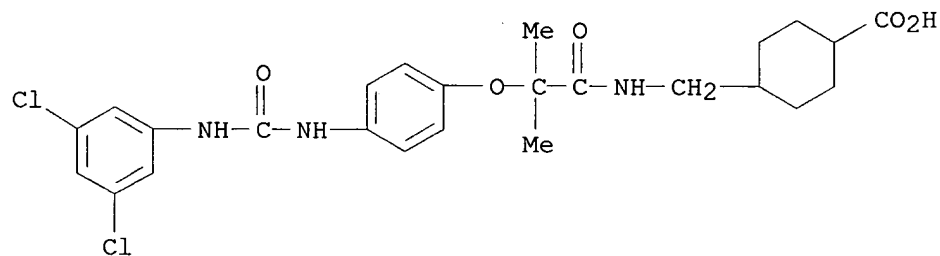


PAGE 1-B



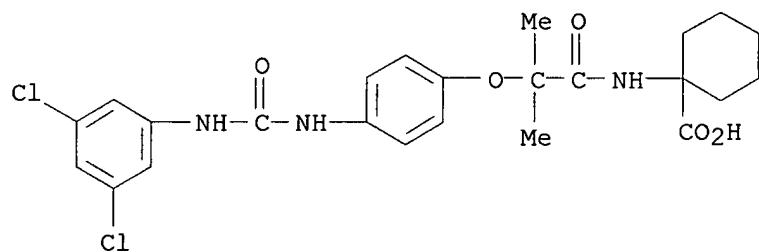
RN 249513-67-5 CAPLUS

CN Cyclohexanecarboxylic acid, 4-[[[2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl-1-oxopropyl]amino]methyl]- (9CI) (CA INDEX NAME)



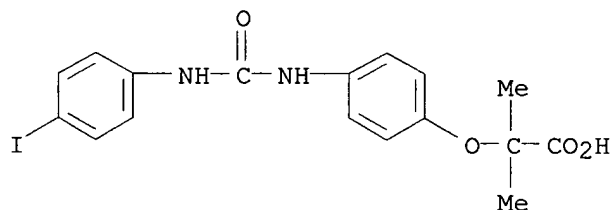
RN 249513-69-7 CAPLUS

CN Cyclohexanecarboxylic acid, 1-[[2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl-1-oxopropyl]amino]- (9CI) (CA INDEX NAME)



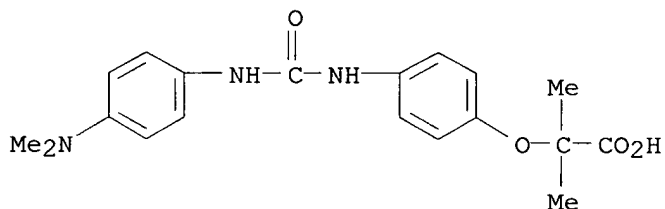
RN 249513-70-0 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-iodophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



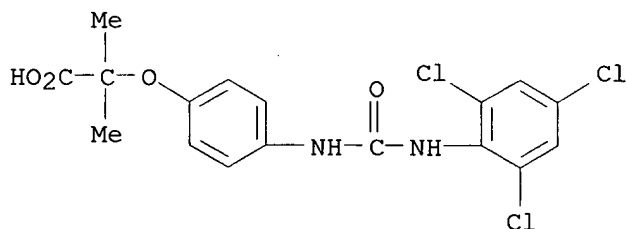
RN 249513-71-1 CAPLUS

CN Propanoic acid, 2-[4-[[[4-(dimethylamino)phenyl]amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



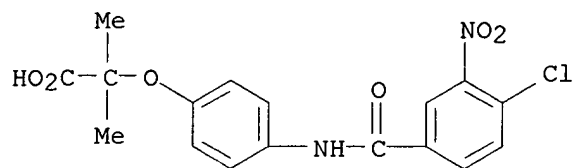
RN 249513-72-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,4,6-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



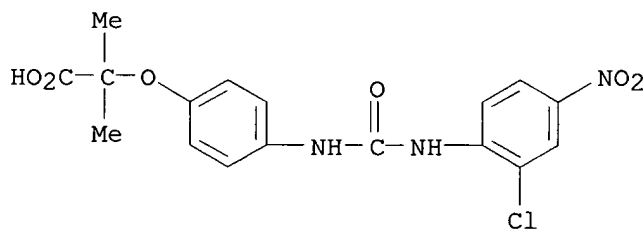
RN 249513-73-3 CAPLUS

CN Propanoic acid, 2-[4-[(4-chloro-3-nitrobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



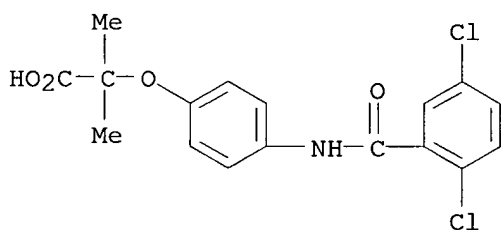
RN 249513-74-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-chloro-4-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



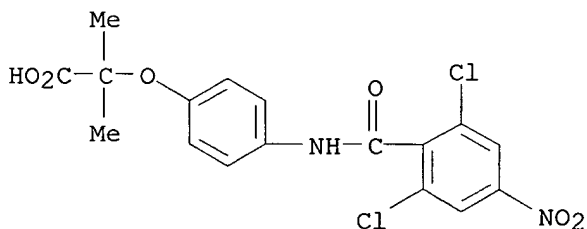
RN 249513-76-6 CAPLUS

CN Propanoic acid, 2-[4-[(2,5-dichlorobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



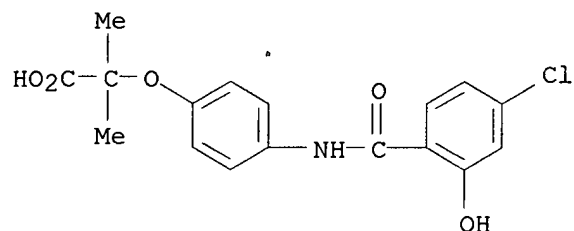
RN 249513-77-7 CAPLUS

CN Propanoic acid, 2-[4-[(2,6-dichloro-4-nitrobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



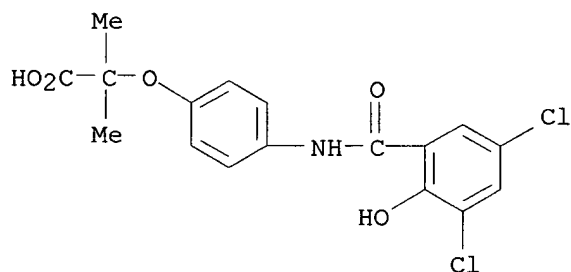
RN 249513-80-2 CAPLUS

CN Propanoic acid, 2-[4-[(4-chloro-2-hydroxybenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



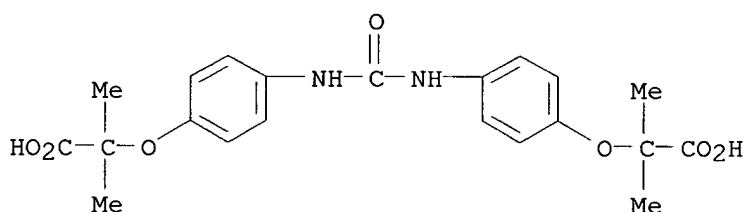
RN 249513-81-3 CAPLUS

CN Propanoic acid, 2-[4-[(3,5-dichloro-2-hydroxybenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



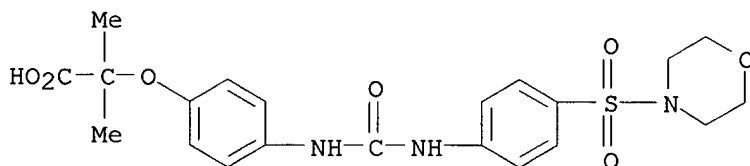
RN 249513-83-5 CAPLUS

CN Propanoic acid, 2,2'-[carbonylbis(imino-4,1-phenyleneoxy)]bis[2-methyl- (9CI) (CA INDEX NAME)



RN 249513-84-6 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[4-(4-morpholinylsulfonyl)phenyl]amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



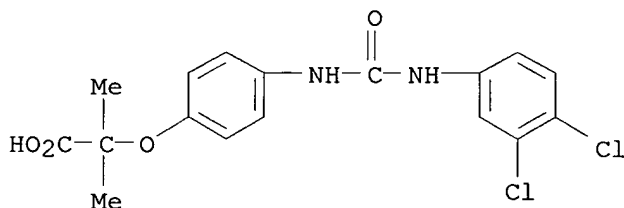
RE.CNT 48

RE

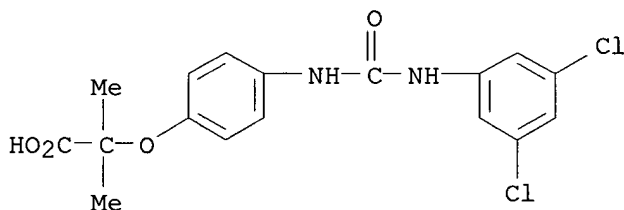
(1) Al-Abed, Y; J Biol Chem 1996, V271, P2892 CAPLUS

- (4) Baynes, J; Diabetes 1999, V48, P1 CAPLUS
 (5) Beisswenger, P; Diabetes 1999, V48, P198 CAPLUS
 (7) Booth, A; J Biol Chem 1997, V272, P5430 CAPLUS
 (8) Brownlee, M; Science 1986, V232, P1629 CAPLUS
 ALL CITATIONS AVAILABLE IN THE RE FORMAT

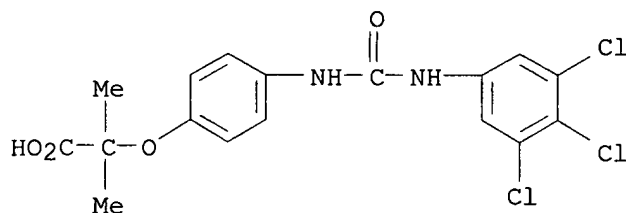
L6 ANSWER 5 OF 19 CAPLUS COPYRIGHT 2000 ACS
 AN 1999:19928 CAPLUS
 DN 130:217602
 TI Structure/function of allosteric effectors of hemoglobin
 AU Abraham, Donald J.; Kister, Jean; Joshi, Gajanan S.; Wireko, Fred C.;
 Safo, Martin K.; Marden, Michael C.; Poyart, Claude
 CS Institute for Structural Biology and Drug Discovery, and Department of
 Medicinal Chemistry, Virginia Commonwealth University, Richmond, VA,
 23298, USA
 SO Med. Chem. Res. (1998), 8(7/8), 478-486
 CODEN: MCREEB; ISSN: 1054-2523
 PB Birkhaeuser Boston
 DT Journal
 LA English
 IT 117011-50-4 121809-80-1, L 35 121809-82-3
 RL: BAC (Biological activity or effector, except adverse); PRP
 (Properties); BIOL (Biological study)
 (structure/function of allosteric effectors of Hb)
 RN 117011-50-4 CAPLUS
 CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
 2-methyl- (9CI) (CA INDEX NAME)



RN 121809-80-1 CAPLUS
 CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
 2-methyl- (9CI) (CA INDEX NAME)



RN 121809-82-3 CAPLUS
 CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trichlorophenyl)amino]carbonyl]ami
 no]phenoxy]- (9CI) (CA INDEX NAME)



RE.CNT 19

RE

- (1) Abraham, D; Biochem 1992, V31, P9141 CAPLUS
- (2) Abraham, D; Blood 1991, V77, P1334 CAPLUS
- (3) Abraham, D; J Mol Biol 1995, V248, P845 CAPLUS
- (4) Abraham, D; Proc Natl Acad Sci 1983, V80, P324 CAPLUS
- (6) Arnone, A; Nature 1972, V237, P146 CAPLUS

ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 6 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1998:207267 CAPLUS

DN 128:270438

TI Preparation of phenoxypropionic acids and analogs as hemoglobin allosteric modifiers

IN Abraham, Donald J.; Joshi, Gajanan; Randad, Ramnarayan; Panikker, Jayashree

PA Virginia Commonwealth University, USA

SO U.S., 43 pp. Cont.-in-part of U.S. 5,432,191.

CODEN: USXXAM

DT Patent

LA English

FAN.CNT 8

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 5731454	A	19980324	US 1995-374206	19950118
	US 5049695	A	19910917	US 1990-478848	19900212
	US 5122539	A	19920616	US 1991-702947	19910520
	US 5382680	A	19950117	US 1991-722382	19910626
	US 5290803	A	19940301	US 1993-6246	19930119
	US 5432191	A	19950711	US 1993-101501	19930730
	US 5591892	A	19970107	US 1995-451658	19950530
	US 5648375	A	19970715	US 1995-478372	19950607
	US 5677330	A	19971014	US 1995-478371	19950607
	US 5705521	A	19980106	US 1995-482808	19950607
	US 5927283	A	19990727	US 1997-848485	19970508
	US 5872282	A	19990216	US 1998-41595	19980313
	PRAI	US 1990-623346		19901207	
US 1990-478848			19901212		
US 1991-702947			19910520		
US 1991-722382			19910626		
US 1993-6246			19930119		
US 1993-101501			19930730		
US 1992-885721			19920518		
US 1995-374206			19950118		
US 1995-478371			19950607		

OS MARPAT 128:270438

IT 121809-80-1P

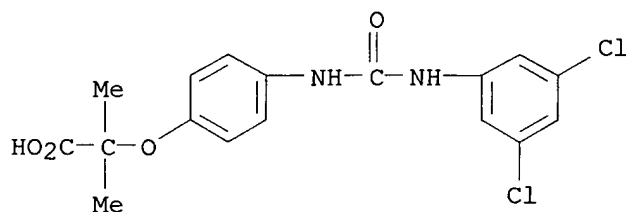
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic

V. Balasubramanian

preparation); THU (Therapeutic use); BIOL (Biological study); PREP
(Preparation); USES (Uses)
(prepn. of phenoxypropionic acids and analogs as Hb allosteric
modifiers)

RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 7 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1998:41719 CAPLUS

DN 128:110865

TI Use of allosteric hemoglobin modifiers in combination with radiation
therapy to treat carcinogenic tumors

IN Abraham, Donald J.

PA Center for Innovative Technology, USA

SO U.S., 47 pp. Cont.-in-part of U.S. Ser. No. 374,206.

CODEN: USXXAM

DT Patent

LA English

FAN.CNT 8

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 5705521	A	19980106	US 1995-482808	19950607
	US 5049695	A	19910917	US 1990-478848	19900212
	US 5122539	A	19920616	US 1991-702947	19910520
	US 5382680	A	19950117	US 1991-722382	19910626
	US 5290803	A	19940301	US 1993-6246	19930119
	US 5432191	A	19950711	US 1993-101501	19930730
	US 5731454	A	19980324	US 1995-374206	19950118
PRAI	US 1990-478848		19900212		
	US 1990-623346		19901207		
	US 1991-702947		19910520		
	US 1991-722382		19910626		
	US 1993-6246		19930119		
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	US 1995-374206		19950118		
	US 1992-885721		19920518		

OS MARPAT 128:110865

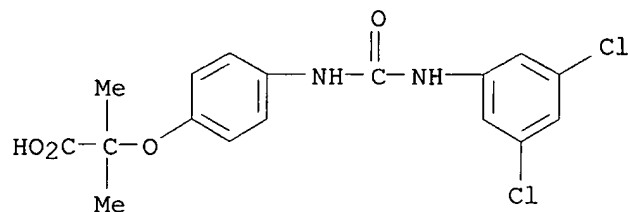
IT **121809-80-1**

RL: BAC (Biological activity or effector, except adverse); THU
(Therapeutic use); BIOL (Biological study); USES (Uses)

(allosteric Hb modifiers in combination with radiation therapy to treat
carcinogenic tumors, prepn., and use for other oxygen tension-related
diseases)

RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 8 OF 19 CAPLUS COPYRIGHT 2000 ACS
 AN 1997:693601 CAPLUS
 DN 127:336654
 TI Medical uses of allosteric hemoglobin modifier compounds in patient care
 IN Abraham, Donald J.; Gerber, Michael
 PA Center for Innovative Technology, USA; Allos Therapeutics, Inc.
 SO U.S., 48 pp. Cont.-in-part of U.S. 5,382,680.
 CODEN: USXXAM
 DT Patent
 LA English
 FAN.CNT 8

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 5677330	A	19971014	US 1995-478371	19950607
	US 5049695	A	19910917	US 1990-478848	19900212
	US 5122539	A	19920616	US 1991-702947	19910520
	US 5382680	A	19950117	US 1991-722382	19910626
	US 5290803	A	19940301	US 1993-6246	19930119
	US 5432191	A	19950711	US 1993-101501	19930730
	US 5731454	A	19980324	US 1995-374206	19950118
	US 5927283	A	19990727	US 1997-848485	19970508
PRAI	US 1990-478848		19900212		
	US 1990-623346		19901207		
	US 1991-702947		19910520		
	US 1991-722382		19910626		
	US 1993-6246		19930119		
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	US 1995-374206		19950118		
	US 1992-885721		19920518		
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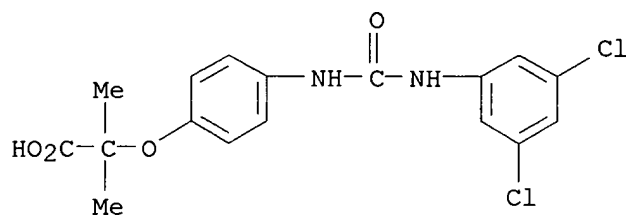
OS MARPAT 127:336654

IT **121809-80-1P**

RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)
 (allosteric Hb modifier compds.)

RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 9 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1994:579000 CAPLUS

DN 121:179000

TI 13C NMR spectra of allosteric effectors of hemoglobin

AU Ostrowski, Stanislaw; Priebe, Waldemar; Burke, Thomas G.

CS M. D. Anderson Cancer Cent., Univ. Texas, Houston, TX, 77030, USA

SO Magn. Reson. Chem. (1994), 32(3), 182-3

CODEN: MRCHEG; ISSN: 0749-1581

DT Journal

LA English

IT 117011-50-4 121809-75-4 121809-83-4

157579-93-6 157579-95-8 157579-96-9

157579-99-2 157580-01-3 157580-02-4

157580-04-6 157580-05-7 157580-07-9

157580-09-1 157580-14-8 157580-16-0

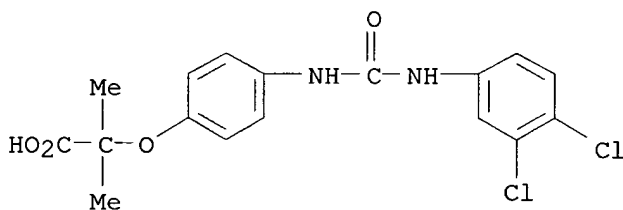
157580-19-3

RL: PRP (Properties)

(NMR of carbon-13 of)

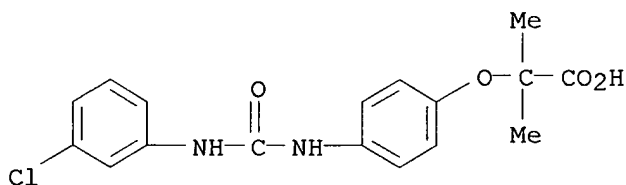
RN 117011-50-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-75-4 CAPLUS

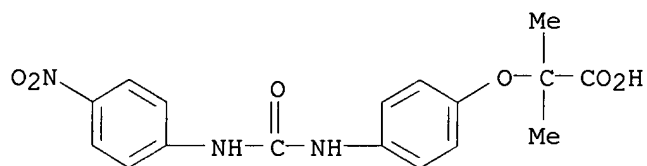
CN Propanoic acid, 2-[4-[[[(3-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-83-4 CAPLUS

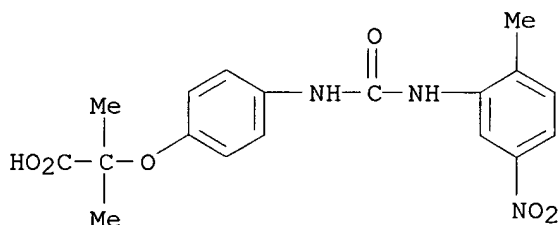
CN Propanoic acid, 2-methyl-2-[4-[[[(4-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

xy]- (9CI) (CA INDEX NAME)



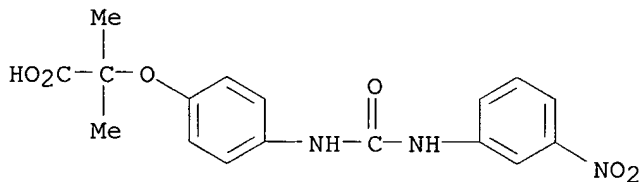
RN 157579-93-6 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2-methyl-5-nitrophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



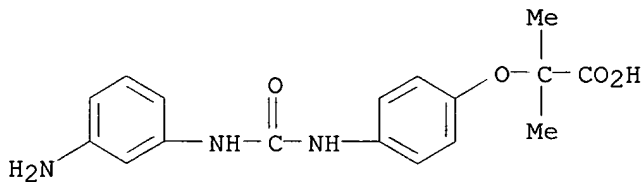
RN 157579-95-8 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3-nitrophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



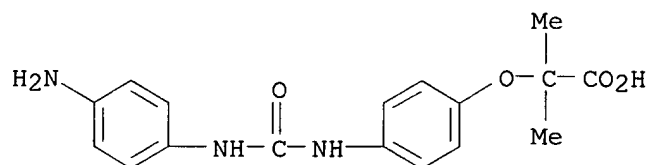
RN 157579-96-9 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-aminophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



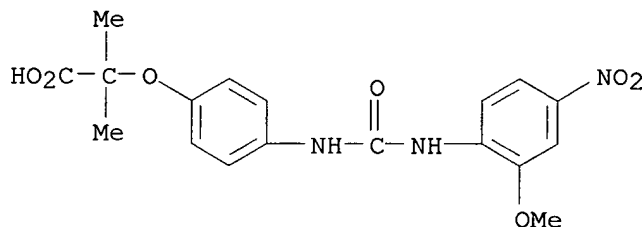
RN 157579-99-2 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-aminophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



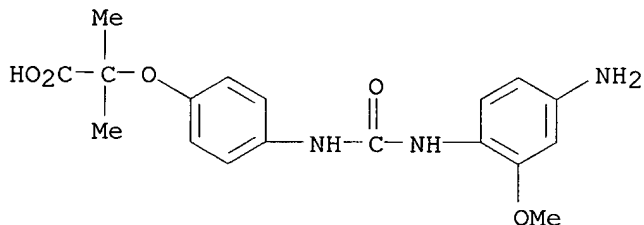
RN 157580-01-3 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-methoxy-4-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



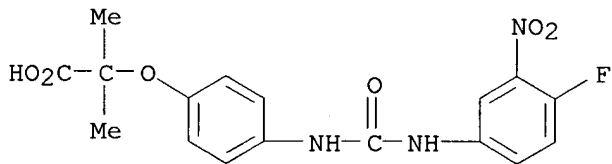
RN 157580-02-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-amino-2-methoxyphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



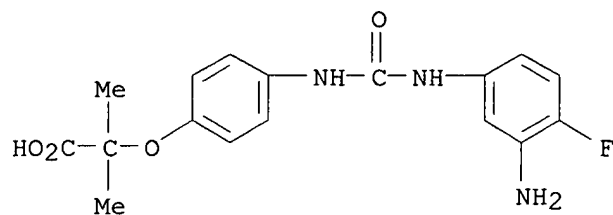
RN 157580-04-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-fluoro-3-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



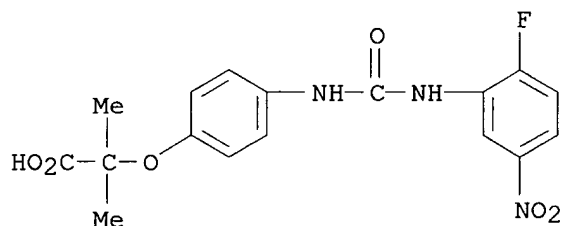
RN 157580-05-7 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-amino-4-fluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



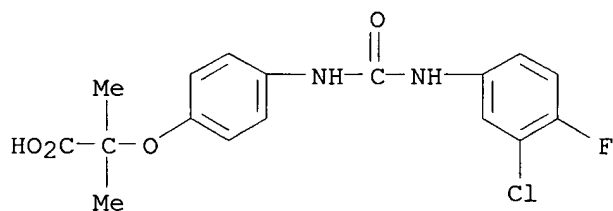
RN 157580-07-9 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-fluoro-5-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



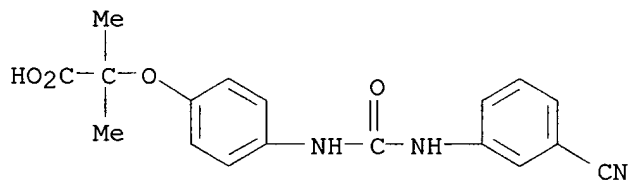
RN 157580-09-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-chloro-4-fluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



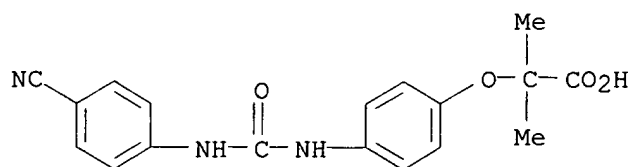
RN 157580-14-8 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-cyanophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



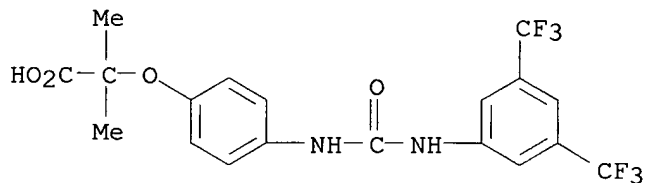
RN 157580-16-0 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-cyanophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 157580-19-3 CAPLUS

CN Propanoic acid, 2-[4-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 10 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1993:455833 CAPLUS

DN 119:55833

TI Evaluation of LEH/LR16 artificial blood formulation

AU Burke, T. G.; Staubus, A. E.

CS Coll. Pharm., Ohio State Univ., Columbus, OH, USA

SO Report (1992), Order No. AD-A248575, 10 pp. Avail.: NTIS

From: Gov. Rep. Announce. Index (U. S.) 1992, 92(15), Abstr. No. 240,053

DT Report

LA English

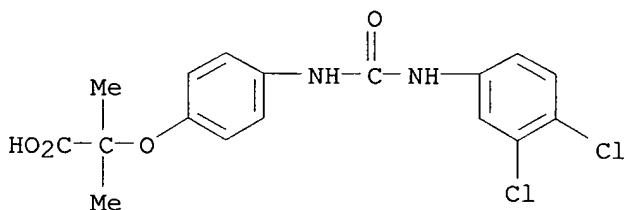
IT **117011-50-4D**, LR 16, analogs

RL: BIOL (Biological study)

(liposomes contg. Hbs and, for blood substitutes)

RN 117011-50-4 CAPLUS

CN Propanoic acid, 2-[4-[[[3,4-dichlorophenyl]amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 11 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1991:639430 CAPLUS

DN 115:239430

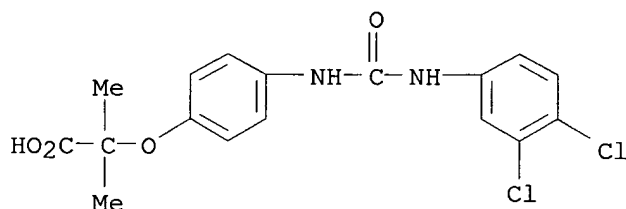
TI Evaluation of liposome-encapsulated hemoglobin/LR16 formulation as a potential blood substitute

AU Burke, T. G.

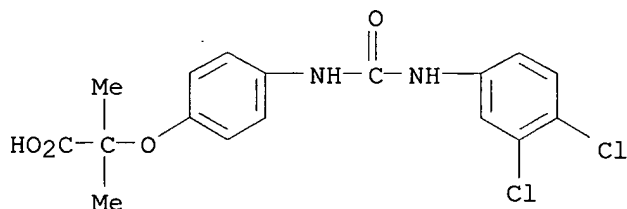
CS City of Hope Natl. Med. Cent., Duarte, CA, USA

V. Balasubramanian

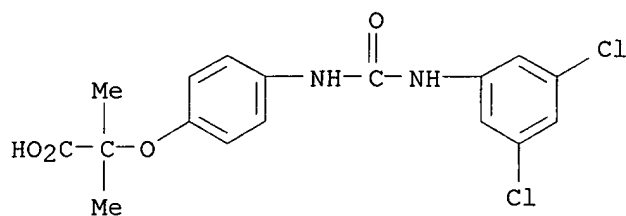
SO Report (1990), Order No. AD-A229 090, 9 pp. Avail.: NTIS
From: Gov. Rep. Announce. Index (U. S.) 1991, 91(10), Abstr. No. 124,831
DT Report
LA English
IT **117011-50-4**, LR 16
RL: BIOL (Biological study)
(liposomes contg. Hb and, for blood substitutes)
RN 117011-50-4 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 12 OF 19 CAPLUS COPYRIGHT 2000 ACS
AN 1991:77296 CAPLUS
DN 114:77296
TI Allosteric modifiers of hemoglobin. 2. Crystallographically determined binding sites and hydrophobic binding/interaction analysis of novel hemoglobin oxygen effectors
AU Wireko, Fred C.; Kellogg, Glen E.; Abraham, Donald J.
CS Med. Coll. Virginia, Virginia Commonw. Univ., Richmond, VA, 23298-0581, USA
SO J. Med. Chem. (1991), 34(2), 758-67
CODEN: JMCMAR; ISSN: 0022-2623
DT Journal
LA English
IT **117011-50-4 121809-80-1 121809-82-3**
RL: BIOL (Biological study)
(Hb of human binding by, hydrophobicity in and oxygen binding response to and allosteric modifier structure relation to)
RN 117011-50-4 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)

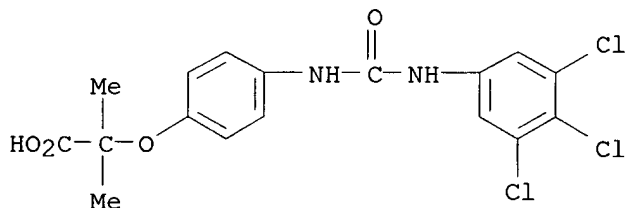


RN 121809-80-1 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



RN 121809-82-3 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



L6 ANSWER 13 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1990:536008 CAPLUS

DN 113:136008

TI Effect of nonionic surfactants on chalcopryrite leaching under dump chemical conditions

AU Sandoval, S. P.; Pool, D. L.; Schultze, L. E.

CS Reno Res. Cent., U. S. Bur. Mines, Reno, NV, USA

SO Bur. Mines Rep. Invest. (1990), RI 9311, 16 pp.

CODEN: XBMIA6; ISSN: 0096-1922

DT Report

LA English

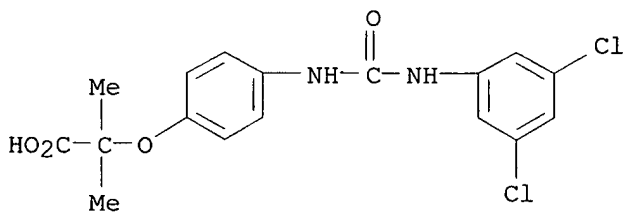
IT 121809-80-1, L35

RL: PROC (Process)

(surfactant, leaching of copper from chalcopryrite with ferric sulfate soln. contg. nonionic)

RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 14 OF 19 CAPLUS COPYRIGHT 2000 ACS

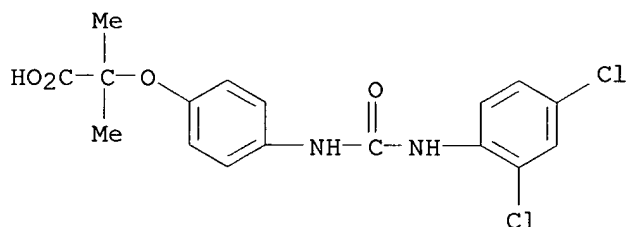
AN 1990:423397 CAPLUS

DN 113:23397

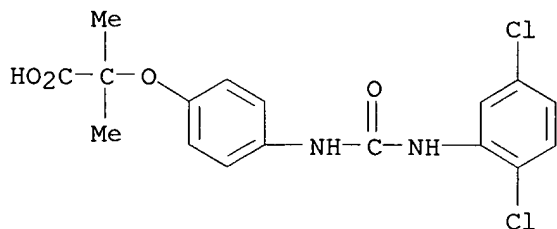
V. Balasubramanian

TI Preparation of (aryluroidphenoxy)- and (benzamidophenoxy)methylpropionic
acid derivatives which reduce the affinity of hemoglobin for oxygen
IN Lalezari, Iraj; Lalezari, Parviz
PA USA
SO PCT Int. Appl., 32 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 8912622	A1	19891228	WO 1989-US2621	19890615
	W: JP				
	RW: AT, BE, CH, DE, FR, GB, IT, LU, NL, SE				
	US 4921997	A	19900501	US 1988-207098	19880615
	EP 420930	A1	19910410	EP 1989-907939	19890615
	R: AT, BE, CH, DE, FR, GB, IT, LI, LU, NL, SE				
	JP 03505728	T2	19911212	JP 1989-507363	19890615
	US 5093367	A	19920303	US 1990-477048	19900207
	US 5292935	A	19940308	US 1992-821409	19920115
	US 5472981	A	19951205	US 1993-158702	19931129
PRAI	US 1988-207098	19880615			
	WO 1989-US2621	19890615			
	US 1990-477048	19900207			
	US 1992-821409	19920115			
OS	MARPAT 113:23397				
IT	121809-78-7 121809-79-8				
	RL: RCT (Reactant)				
	(Hb-oxygen affinity, effect on)				
RN	121809-78-7 CAPLUS				
CN	Propanoic acid, 2-[4-[[[(2,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]- 2-methyl- (9CI) (CA INDEX NAME)				



RN 121809-79-8 CAPLUS
CN Propanoic acid, 2-[4-[[[(2,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)

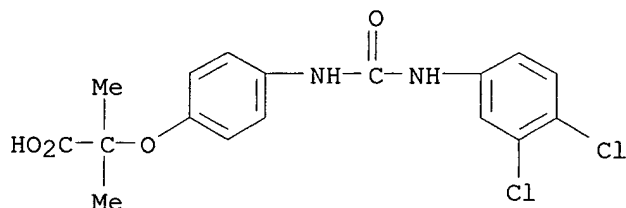


IT 117011-50-4P 121809-73-2P 121809-80-1P
121809-81-2P 121809-82-3P 121809-91-4P
127676-57-7P

RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. of, as Hb-oxygen affinity-reducing agent)

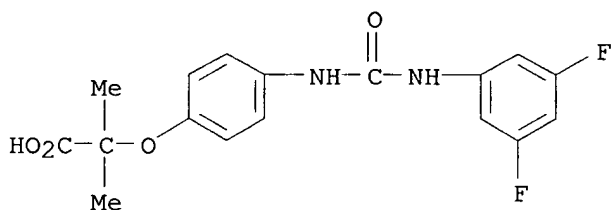
RN 117011-50-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



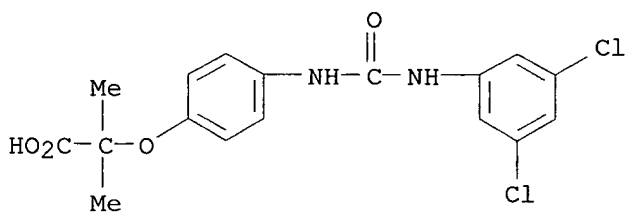
RN 121809-73-2 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-difluorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



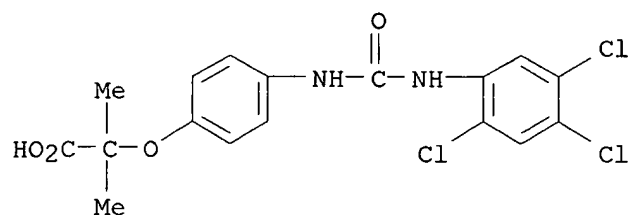
RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



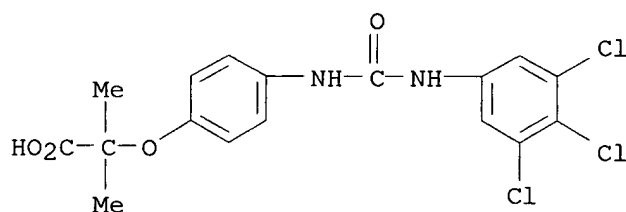
RN 121809-81-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



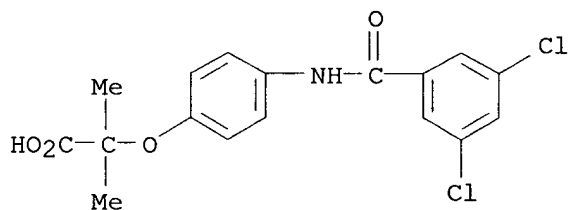
RN 121809-82-3 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



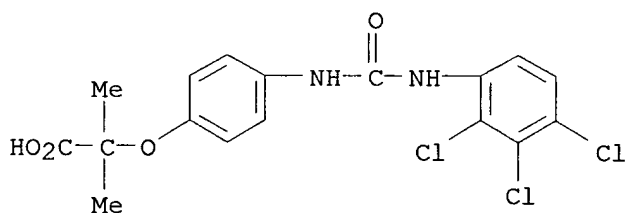
RN 121809-91-4 CAPLUS

CN Propanoic acid, 2-[4-[(3,5-dichlorobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 127676-57-7 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,3,4-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



L6 ANSWER 15 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1990:231450 CAPLUS

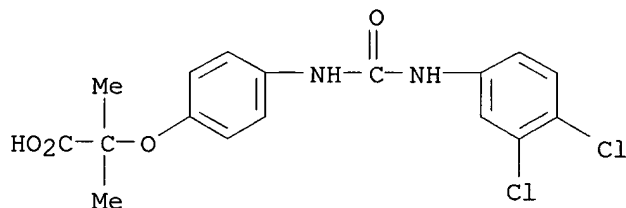
DN 112:231450

TI Effectors of hemoglobin. Separation of allosteric and affinity factors

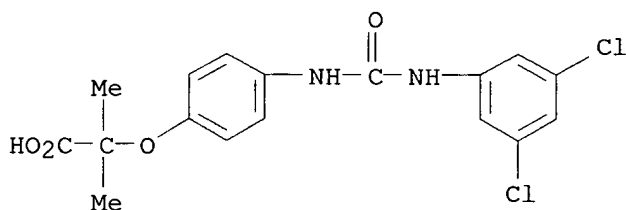
AU Marden, Michael C.; Bohn, Brigitte; Kister, Jean; Poyart, Claude

V. Balasubramanian

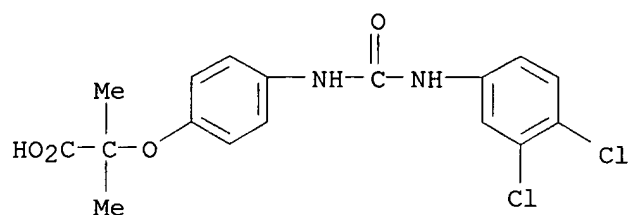
CS Hop. Bicetre, Le Kremlin Bicetre, 94275, Fr.
SO Biophys. J. (1990), 57(3), 397-403
CODEN: BIOJAU; ISSN: 0006-3495
DT Journal
LA English
IT **117011-50-4 121809-80-1**, L 35
RL: BIOL (Biological study)
(Hb binding by carbon monoxide and oxygen response to, effectors
affinity and allosterism in)
RN 117011-50-4 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



RN 121809-80-1 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



L6 ANSWER 16 OF 19 CAPLUS COPYRIGHT 2000 ACS
AN 1990:72472 CAPLUS
DN 112:72472
TI New effectors of human hemoglobin: structure and function
AU Lalezari, I.; Lalezari, P.; Poyart, C.; Marden, M.; Kister, J.; Bohn, B.;
Fermi, G.; Perutz, M. F.
CS Montefiore Med. Cent., Albert Einstein Coll. Med., Bronx, NY, 10467, USA
SO Biochemistry (1990), 29(6), 1515-23
CODEN: BICHAW; ISSN: 0006-2960
DT Journal
LA English
IT **117011-50-4**, LR16
RL: BIOL (Biological study)
(as Hb A of human allosteric effector, methylpropionic acid derivs.
compared to)
RN 117011-50-4 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



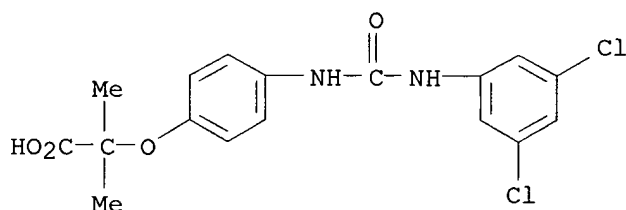
IT 121809-80-1 121809-82-3

RL: PROC (Process)

(as Hb A of human allosteric effector, structure and function of)

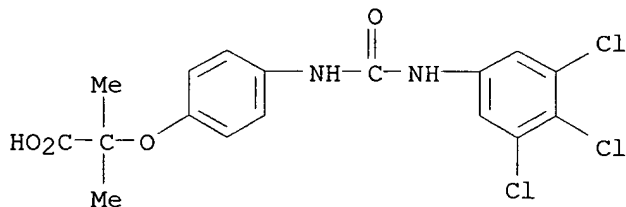
RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-82-3 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



L6 ANSWER 17 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1989:587608 CAPLUS

DN 111:187608

TI Compound, composition, and method for the reduction of lipids, the modification of the affinity of hemoglobin for oxygen, and the prevention of platelet aggregation

IN Lalezari, Iraj; Rahbar, Samuel; Lalezari, Parviz

PA USA

SO PCT Int. Appl., 15 pp.

CODEN: PIXXD2

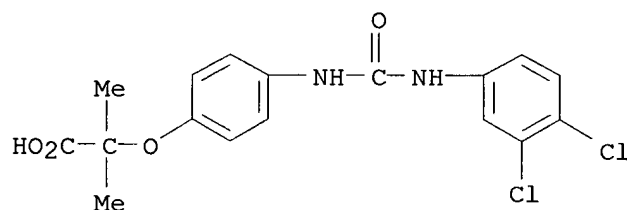
DT Patent

LA English

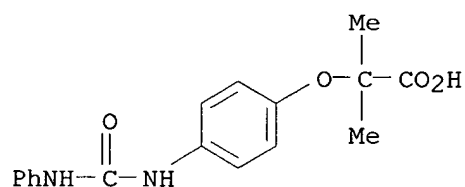
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 8810113	A1	19881229	WO 1988-US2092	19880615
	W: AU, JP				

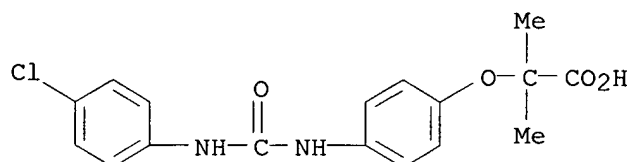
RW: AT, BE, CH, DE, FR, GB, IT, LU, NL, SE
 AU 8819986 A1 19890119 AU 1988-19986 19880615
 AU 616674 B2 19911107
 EP 358724 A1 19900321 EP 1988-906435 19880615
 R: AT, BE, CH, DE, FR, GB, IT, LI, LU, NL, SE
 JP 02503797 T2 19901108 JP 1988-505967 19880615
 US 5268500 A 19931207 US 1991-778706 19911018
 PRAI US 1987-62236 19870615
 WO 1988-US2092 19880615
 US 1989-327020 19890322
 US 1990-515673 19900412
 US 1991-658096 19910220
 OS MARPAT 111:187608
 IT **117011-50-4P 121809-57-2P 121809-76-5P**
 RL: SPN (Synthetic preparation); PREP (Preparation)
 (prepn. of, as antihyperlipidemic and for modification of oxygen
 affinity of Hb)
 RN 117011-50-4 CAPLUS
 CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
 2-methyl- (9CI) (CA INDEX NAME)



RN 121809-57-2 CAPLUS
 CN Propanoic acid, 2-methyl-2-[4-[(phenylamino)carbonyl]amino]phenoxy]-
 (9CI) (CA INDEX NAME)



RN 121809-76-5 CAPLUS
 CN Propanoic acid, 2-[4-[[[(4-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-
 methyl- (9CI) (CA INDEX NAME)



V. Balasubramanian

L6 ANSWER 18 OF 19 CAPLUS COPYRIGHT 2000 ACS

AN 1989:533730 CAPLUS

DN 111:133730

TI Synthesis and investigation of effects of 2-[4-
[[(arylamino)carbonyl]amino]phenoxy]-2-methylpropionic acids on the
affinity of hemoglobin for oxygen: structure-activity relationships

AU Lalezari, Iraj; Lalezari, Parviz

CS Dep. Med., Montefiore Med. Cent., Bronx, NY, 10467, USA

SO J. Med. Chem. (1989), 32(10), 2352-7

CODEN: JMCMAR; ISSN: 0022-2623

DT Journal

LA English

OS CASREACT 111:133730

IT 117011-50-4 121809-77-6 121809-78-7

121809-79-8 121809-80-1 121809-81-2

121809-82-3 121809-83-4 121809-84-5

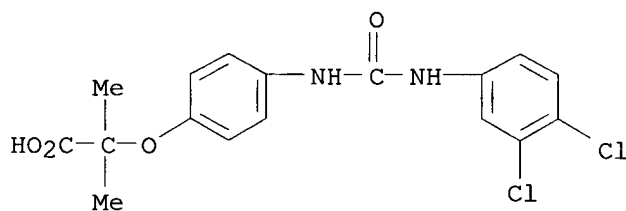
121809-85-6 121809-86-7 121809-87-8

RL: RCT (Reactant)

(prepn. and Hb-oxygen affinity in presence of)

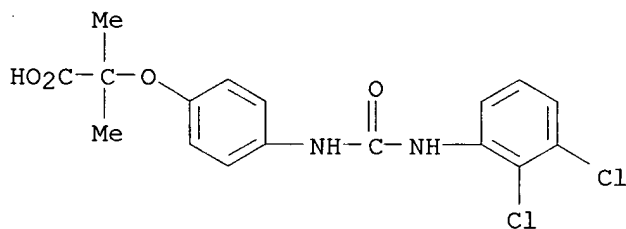
RN 117011-50-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



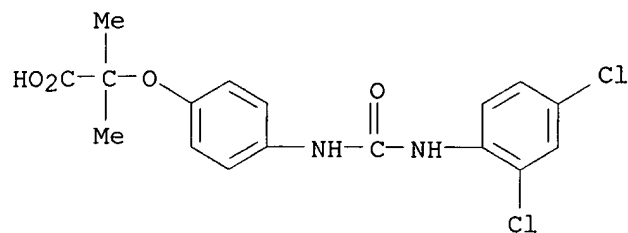
RN 121809-77-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(2,3-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



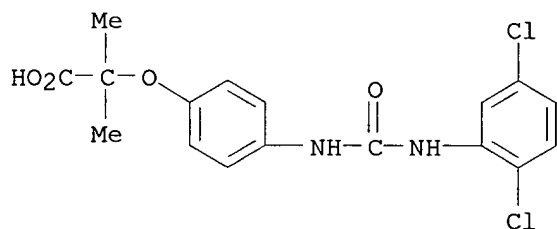
RN 121809-78-7 CAPLUS

CN Propanoic acid, 2-[4-[[[(2,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



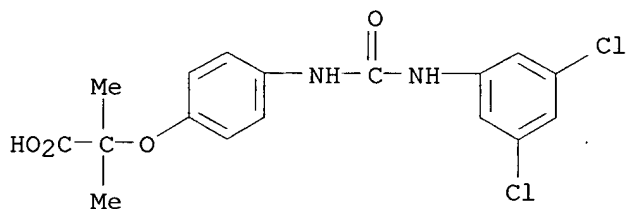
RN 121809-79-8 CAPLUS

CN Propanoic acid, 2-[4-[[[(2,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



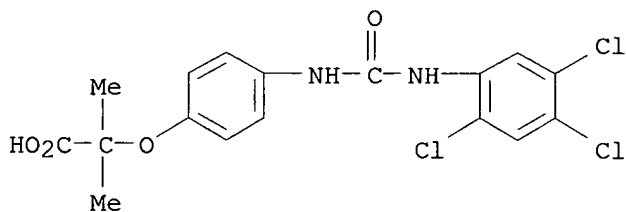
RN 121809-80-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-dichlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



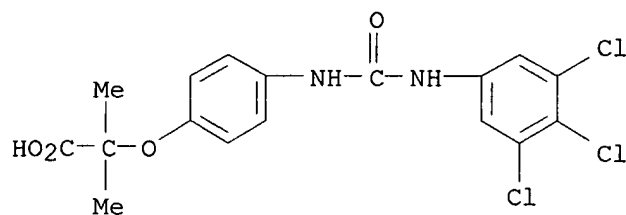
RN 121809-81-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



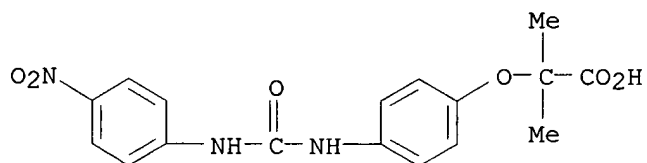
RN 121809-82-3 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trichlorophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



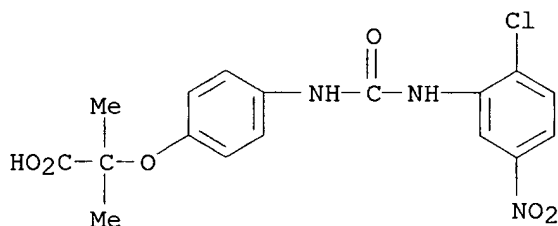
RN 121809-83-4 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(4-nitrophenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



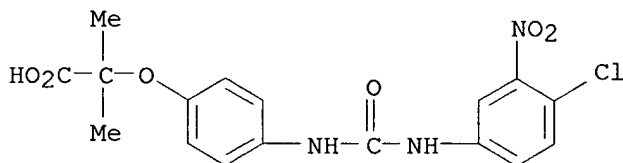
RN 121809-84-5 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-chloro-5-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



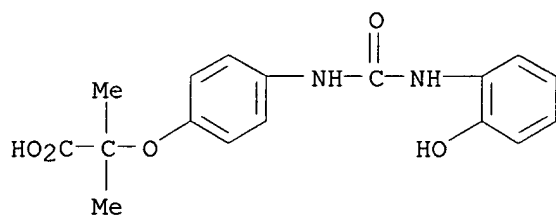
RN 121809-85-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-chloro-3-nitrophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



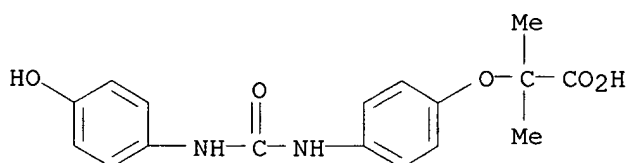
RN 121809-86-7 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-hydroxyphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



RN 121809-87-8 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-hydroxyphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

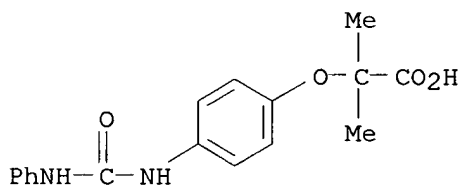


IT 121809-57-2P 121809-58-3P 121809-65-2P
 121809-66-3P 121809-67-4P 121809-68-5P
 121809-69-6P 121809-70-9P 121809-71-0P
 121809-72-1P 121809-73-2P 121809-74-3P
 121809-75-4P 121809-76-5P 121809-88-9P
 121809-89-0P 121809-90-3P 121809-91-4P
 121809-94-7P

RL: SPN (Synthetic preparation); PREP (Preparation)
 (prepn. and Hb-oxygen affinity in the presence of)

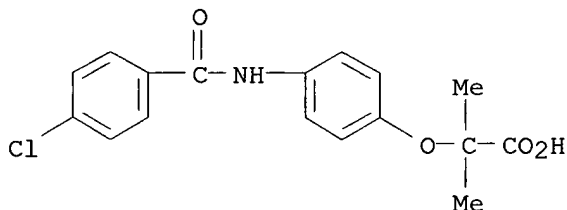
RN 121809-57-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[(phenylamino)carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



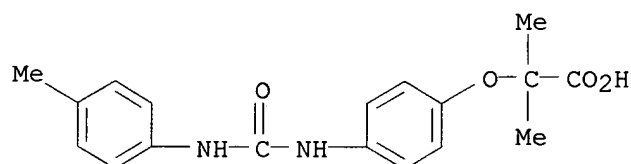
RN 121809-58-3 CAPLUS

CN Propanoic acid, 2-[4-[(4-chlorobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



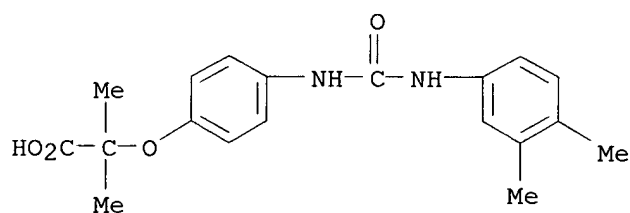
RN 121809-65-2 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(4-methylphenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



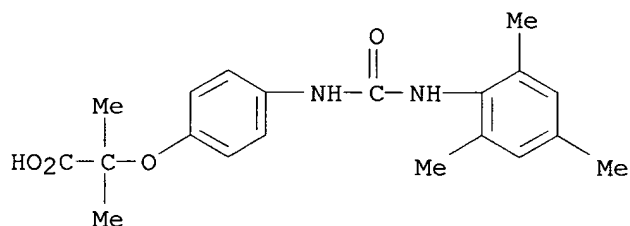
RN 121809-66-3 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,4-dimethylphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



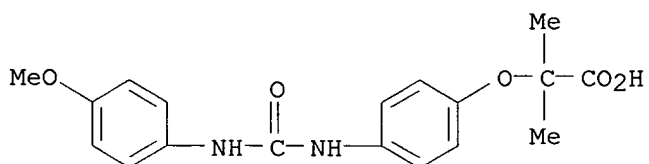
RN 121809-67-4 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(2,4,6-trimethylphenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



RN 121809-68-5 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-methoxyphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

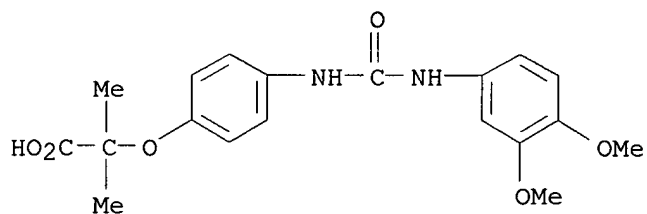


RN 121809-69-6 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,4-dimethoxyphenyl)amino]carbonyl]amino]phenoxy]-

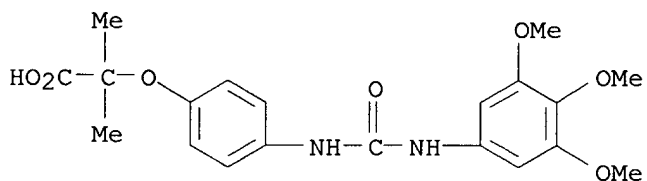
V. Balasubramanian

2-methyl- (9CI) (CA INDEX NAME)



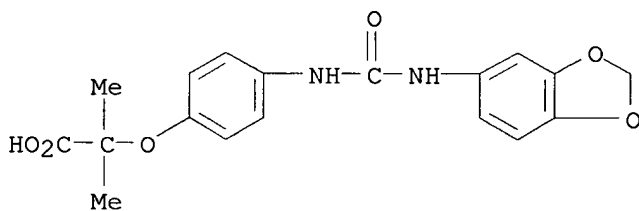
RN 121809-70-9 CAPLUS

CN Propanoic acid, 2-methyl-2-[4-[[[(3,4,5-trimethoxyphenyl)amino]carbonyl]amino]phenoxy]- (9CI) (CA INDEX NAME)



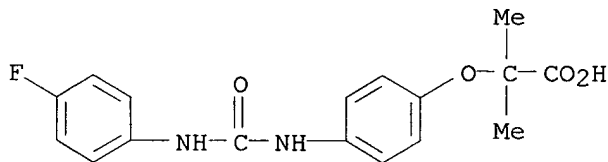
RN 121809-71-0 CAPLUS

CN Propanoic acid, 2-[4-[[[(1,3-benzodioxol-5-ylamino)carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



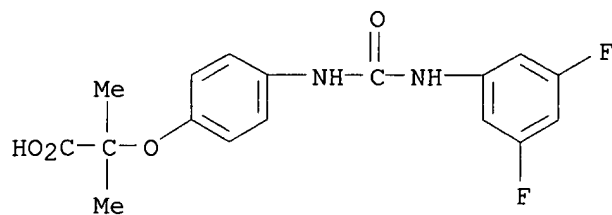
RN 121809-72-1 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-fluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



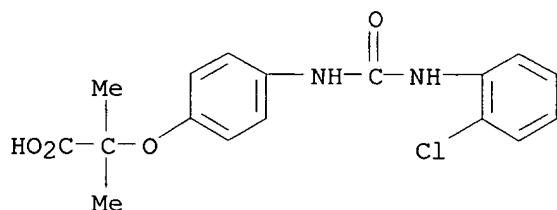
RN 121809-73-2 CAPLUS

CN Propanoic acid, 2-[4-[[[(3,5-difluorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



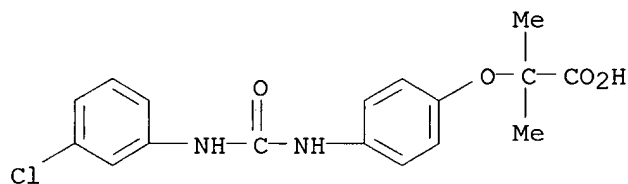
RN 121809-74-3 CAPLUS

CN Propanoic acid, 2-[4-[[[(2-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



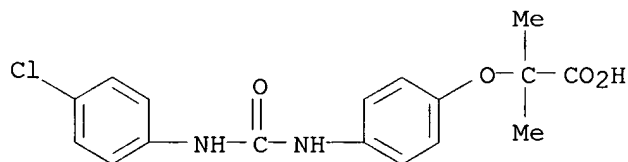
RN 121809-75-4 CAPLUS

CN Propanoic acid, 2-[4-[[[(3-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



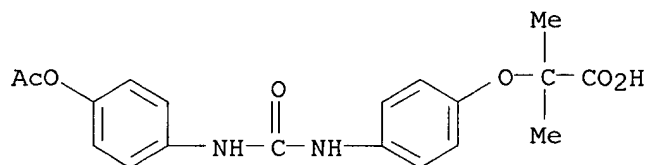
RN 121809-76-5 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-chlorophenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



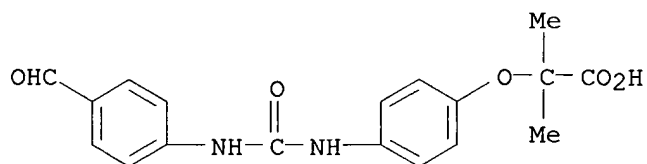
RN 121809-88-9 CAPLUS

CN Propanoic acid, 2-[4-[[[(4-(acetyloxy)phenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



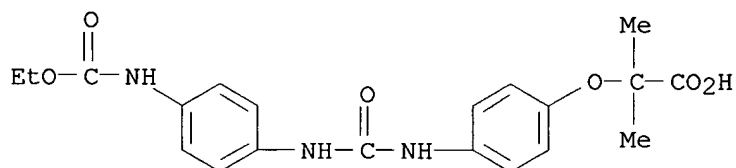
RN 121809-89-0 CAPLUS

CN Propanoic acid, 2-[4-[[[4-(4-formylphenyl)amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



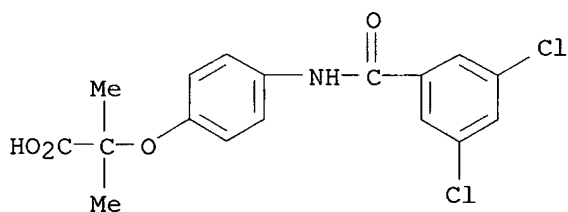
RN 121809-90-3 CAPLUS

CN Propanoic acid, 2-[4-[[[4-[(ethoxycarbonyl)amino]phenyl]amino]carbonyl]amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)



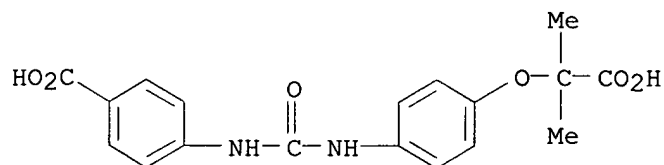
RN 121809-91-4 CAPLUS

CN Propanoic acid, 2-[4-[(3,5-dichlorobenzoyl)amino]phenoxy]-2-methyl- (9CI) (CA INDEX NAME)

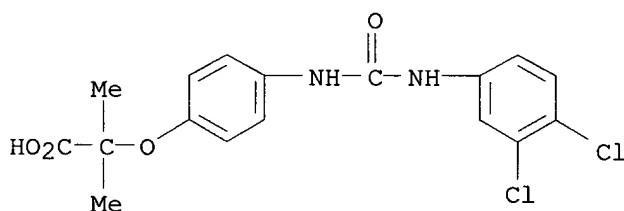


RN 121809-94-7 CAPLUS

CN Benzoic acid, 4-[[[4-(1-carboxy-1-methylethoxy)phenyl]amino]carbonyl]amino]- (9CI) (CA INDEX NAME)



L6 ANSWER 19 OF 19 CAPLUS COPYRIGHT 2000 ACS
AN 1988:563263 CAPLUS
DN 109:163263
TI LR16, a compound with potent effects on the oxygen affinity of hemoglobin,
on blood cholesterol, and on low density lipoprotein
AU Lalezari, I.; Rahbar, S.; Lalezari, P.; Fermi, G.; Perutz, M. F.
CS Dep. Anesthesiol., Montefiore Med. cent., Bronx, NY, 10467, USA
SO Proc. Natl. Acad. Sci. U. S. A. (1988), 85(16), 6117-21
CODEN: PNASA6; ISSN: 0027-8424
DT Journal
LA English
IT **117011-50-4P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. and anticholesterolemic activity of and Hb oxygen affinity
response to)
RN 117011-50-4 CAPLUS
CN Propanoic acid, 2-[4-[[[(3,4-dichlorophenyl)amino]carbonyl]amino]phenoxy]-
2-methyl- (9CI) (CA INDEX NAME)



=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	53.21	179.96

STN INTERNATIONAL LOGOFF AT 15:56:19 ON 05 OCT 2000

Trying 3106016892...Open

Welcome to STN International! Enter x:x
LOGINID:sssptal611bxv
PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

V. Balasubramanian

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 Dec 17 The CA Lexicon available in the CAPLUS and CA files
NEWS 3 Feb 06 Engineering Information Encompass files have new names
NEWS 4 Feb 16 TOXLINE no longer being updated
NEWS 5 Apr 23 Search Derwent WPINDEX by chemical structure
NEWS 6 Apr 23 PRE-1967 REFERENCES NOW SEARCHABLE IN CAPLUS AND CA
NEWS 7 May 07 DGENE Reload
NEWS 8 Jun 20 Published patent applications (A1) are now in USPATFULL
NEWS 9 JUL 13 New SDI alert frequency now available in Derwent's
DWPI and DPCI
NEWS 10 Aug 23 In-process records and more frequent updates now in
MEDLINE
NEWS 11 Aug 23 PAGE IMAGES FOR 1947-1966 RECORDS IN CAPLUS AND CA
NEWS 12 Aug 23 Adis Newsletters (ADISNEWS) now available on STN
NEWS 13 Sep 17 IMSworld Pharmaceutical Company Directory name change
to PHARMASEARCH

NEWS EXPRESS August 15 CURRENT WINDOWS VERSION IS V6.0c,
CURRENT MACINTOSH VERSION IS V6.0 (ENG) AND V6.0J (JP),
AND CURRENT DISCOVER FILE IS DATED 07 AUGUST 2001
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NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
specific topic.

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FILE 'HOME' ENTERED AT 14:20:49 ON 23 SEP 2001

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.15

0.15

FILE 'CAPLUS' ENTERED AT 14:20:58 ON 23 SEP 2001

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26, 1996), unless otherwise indicated in the original publications.

FILE COVERS 1947 - 23 Sep 2001 VOL 135 ISS 14

FILE LAST UPDATED: 21 Sep 2001 (20010921/ED)

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=> s advanced(l)glycation(l)ednproduct?

71854 ADVANCED

2722 GLYCATION

0 EDNPRODUCT?

L1 0 ADVANCED(L) GLYCATION(L) EDNPRODUCT?

=> s advanced(l)glycation(l)endproduct?

71854 ADVANCED

2722 GLYCATION

571 ENDPRODUCT?

L2 233 ADVANCED(L) GLYCATION(L) ENDPRODUCT?

=> l2 and acid

L2 IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system.

For a list of commands available to you in the current file, enter

"HELP COMMANDS" at an arrow prompt (=>).

=> s l2 and acid

3008179 ACID

L3 35 L2 AND ACID

=> d scan l3

L3 35 ANSWERS CAPLUS COPYRIGHT 2001 ACS

CC 14-11 (Mammalian Pathological Biochemistry)

Section cross-reference(s): 2, 15

TI The receptor for advanced glycation end products mediates the chemotaxis of rabbit smooth muscle cells

ST **advanced glycation endproduct** receptor

smooth muscle; AGE receptor vascular smooth muscle atherosclerosis;

TGFbeta AGE receptor smooth muscle atherosclerosis; chemotaxis AGE

receptor smooth muscle atherosclerosis

(lesion; receptor for advanced glycation end products mediates chemotaxis of rabbit smooth muscle cells in relation to receptor-mediated endocytosis, transforming growth factor-.beta. secretion, foam cells, and atherosclerotic lesions)
(receptor for advanced glycation end products mediates chemotaxis of rabbit smooth muscle cells in relation to receptor-mediated endocytosis, transforming growth factor-.beta. secretion, foam cells, and atherosclerotic lesions)
(receptor for advanced glycation end products mediates chemotaxis of rabbit smooth muscle cells in relation to receptor-mediated endocytosis, transforming growth factor-.beta. secretion, foam cells, and atherosclerotic lesions)
(receptors; receptor for advanced glycation end products mediates chemotaxis of rabbit smooth muscle cells in relation to receptor-mediated endocytosis, transforming growth factor-.beta. secretion, foam cells, and atherosclerotic lesions)
(receptors; receptor for advanced glycation end products mediates chemotaxis of rabbit smooth muscle cells in relation to receptor-mediated endocytosis, transforming growth factor-.beta. secretion, foam cells, and atherosclerotic lesions)

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> s 12 and (benzoic(1)acid or phenoxy(1)isobutyric(1)acid)

42974 BENZOIC

3008179 ACID

39392 BENZOIC(L)ACID

18168 PHENOXY

5445 ISOBUTYRIC

3008179 ACID

24 PHENOXY(L)ISOBUTYRIC(L)ACID

L4 1 L2 AND (BENZOIC(L)ACID OR PHENOXY(L)ISOBUTYRIC(L)ACID)

=> d scan

L4 1 ANSWERS CAPLUS COPYRIGHT 2001 ACS

CC 1-12 (Pharmacology)

TI Novel Inhibitors of **Advanced Glycation**

Endproducts (Part II)

ST **advanced glycation endproduct** inhibitor

(AGE-protein; inhibitors of **advanced glycation**

endproducts)

(inhibitors of **advanced glycation**

endproducts)

(inhibitors of **advanced glycation**

endproducts)

(inhibitors of advanced glycation endproducts)

ALL ANSWERS HAVE BEEN SCANNED

=> s 12 and benzoic(1)acid

42974 BENZOIC

3008179 ACID

39392 BENZOIC(L)ACID

L5 1 L2 AND BENZOIC(L)ACID

V. Balasubramanian

=> s l2 and phenoxy
18168 PHENOXY
L6 1 L2 AND PHENOXY

=> d l5 bib abs hitstr

L5 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2001 ACS
AN 2000:835182 CAPLUS
DN 134:247197
TI Novel Inhibitors of **Advanced Glycation Endproducts** (Part II)
AU Rahbar, Samuel; Yerneni, Kiran Kumar; Scott, Stephen; Gonzales, Noe; Lalezari, Iraj
CS Department of Diabetes, Endocrinology, and Metabolism, Gonda (Goldschmied) Diabetes and Genetic Research Center, City of Hope National Medical Center, Duarte, CA, 91010, USA
SO Mol. Cell Biol. Res. Commun. (2000), 3(6), 360-366
CODEN: MCBCFS; ISSN: 1522-4724
PB Academic Press
DT Journal
LA English
AB Enhanced formation and accumulation of **advanced glycation endproducts** (AGEs), have been implicated as a major pathogenesis process leading to diabetic complications, normal aging, atherosclerosis, and Alzheimer's Disease. Several potential drug candidates as AGE inhibitors have been reported recently. The aim of this study was to develop classes of novel inhibitors of **glycation**, AGE formation, and AGE-crosslinking and to investigate their effects through in vitro chem. and immunochem. assays. A total of 92 compds. were designed and synthesized. The first 63 compds. were reported before. Nearly half of the 29 novel inhibitors reported here are **benzoic acid** derivs. and related mols., and found to be potent inhibitors of multistage **glycation**, AGE formation, and AGE-protein crosslinking. All 29 compds. show some degrees of inhibitory activities as detected by the four assay methods, 9 compds. demonstrated high percent inhibition (PI) in all tests, 30 to 40 times stronger than aminoguanidine.
(c) 2000 Academic Press.

RE.CNT 50

RE

(1) Al-Abed, Y; J Biol Chem 1996, V271, P2892 CAPLUS
(3) Anon; US 5514676 CAPLUS
(5) Baynes, J; Diabetes 1999, V48, P1 CAPLUS
(8) Bucala, R; Adv Pharmacol 1992, V23, P1 CAPLUS
(9) Bucala, R; Exp Physiol 1997, V82, P327 CAPLUS
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> d l6 bib abs

L6 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2001 ACS
AN 1999:558280 CAPLUS
DN 131:317717
TI Novel Inhibitors of **Advanced Glycation Endproducts**
AU Rahbar, Samuel; Kumar Yernini, Kiran; Scott, Stephen; Gonzales, Noe; Lalezari, Iraj
CS Department of Diabetes, Endocrinology & Metabolism, City of Hope National Medical Center, Duarte, CA, 91010-0269, USA

09/825,925

V. Balasubramanian

SO Biochem. Biophys. Res. Commun. (1999), 262(3), 651-656

CODEN: BBRCA9; ISSN: 0006-291X

PB Academic Press

DT Journal

LA English

AB Enhanced formation and accumulation of **advanced**

glycation endproducts (AGE's) have been proposed to play a major role in the pathogenesis of diabetic complications, aging, atherosclerosis, and Alzheimer disease leading to progressive and irreversible intermol. protein crosslinkings. This process is accelerated in diabetes and has been postulated to contribute to the development of a range of diabetic complications including nephropathy, retinopathy and neuropathy. Several potential drug candidates as AGE inhibitors have been reported recently. Aminoguanidine is the first drug extensively studied both in vitro and in vivo. The authors have developed a new class of compds. as potent inhibitors of **glycation** and AGE formation. The novel inhibitors reported here are aryl (and heterocyclic) ureido, and aryl (and heterocyclic) carboxamido **phenoxy** isobutyric acids and related mols., which were found by in vitro assay methods to be potent inhibitors of multiple stage of **glycation** and AGE formation.

(c) 1999 Academic Press.

RE.CNT 48

RE

(1) Al-Abed, Y; J Biol Chem 1996, V271, P2892 CAPLUS

(4) Baynes, J; Diabetes 1999, V48, P1 CAPLUS

(5) Beisswenger, P; Diabetes 1999, V48, P198 CAPLUS

(7) Booth, A; J Biol Chem 1997, V272, P5430 CAPLUS

(8) Brownlee, M; Science 1986, V232, P1629 CAPLUS

ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

30.26

30.41

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

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-1.18

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* * * * *

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